

# Package ‘ggmix’

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**Type** Package

**Title** Variable Selection in Linear Mixed Models for SNP Data

**Version** 0.0.2

**Description** Fit penalized multivariable linear mixed models with a single random effect to control for population structure in genetic association studies. The goal is to simultaneously fit many genetic variants at the same time, in order to select markers that are independently associated with the response. Can also handle prior annotation information, for example, rare variants, in the form of variable weights. For more information, see the website below and the accompanying paper: Bhatnagar et al., ``Simultaneous SNP selection and adjustment for population structure in high dimensional prediction models", 2020, <[DOI:10.1371/journal.pgen.1008766](https://doi.org/10.1371/journal.pgen.1008766)>.

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**Encoding** UTF-8

**LazyData** true

**Depends** R (>= 3.4.0)

**Imports** glmnet, methods, stats, MASS, Matrix

**Suggests** RSpectra, popkin, bnpsd, testthat, covr, knitr, rmarkdown

**BugReports** <https://github.com/sahirbhatnagar/ggmix/issues>

**URL** <https://github.com/sahirbhatnagar/ggmix>

**RoxygenNote** 7.1.1

**VignetteBuilder** knitr

**NeedsCompilation** no

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admixed	<i>Simulated Dataset with 1D Geography</i>
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Description

A simulated dataset to show the utility of this package

Usage

admixed

Format

An object of class list of length 21.

Details

The code used to simulate the data is available at <https://github.com/sahirbhatnagar/ggmix/blob/master/data-raw/bnpsd-data.R>. See [gen\\_structured\\_model](#) for more details on the output and how the function used to simulate the data.

**Value**

A list with the following elements

**ytrain** simulated response vector for training set

**ytune** simulated response vector for tuning parameter selection set

**ytest** simulated response vector for test set

**xtrain** simulated design matrix for training set

**xtune** simulated design matrix for tuning parameter selection set

**xtest** simulated design matrix for testing set

**xtrain\_lasso** simulated design matrix for training set for lasso model. This is the same as xtrain, but also includes the nPC principal components

**xtune\_lasso** simulated design matrix for tuning parameter selection set for lasso model. This is the same as xtune, but also includes the nPC principal components

**xtest** simulated design matrix for testing set for lasso model. This is the same as xtest, but also includes the nPC principal components

**causal** character vector of the names of the causal SNPs

**beta** the vector of true regression coefficients

**kin\_train** 2 times the estimated kinship for the training set individuals

**kin\_tune\_train** The covariance matrix between the tuning set and the training set individuals

**kin\_test\_train** The covariance matrix between the test set and training set individuals

**Xkinship** the matrix of SNPs used to estimate the kinship matrix

**not\_causal** character vector of the non-causal SNPs

**PC** the principal components for population structure adjustment

**References**

Ochoa, Alejandro, and John D. Storey. 2016a. "FST And Kinship for Arbitrary Population Structures I: Generalized Definitions." *bioRxiv* doi:10.1101/083915.

Ochoa, Alejandro, and John D. Storey. 2016b. "FST And Kinship for Arbitrary Population Structures II: Method of Moments Estimators." *bioRxiv* doi:10.1101/083923.

**Examples**

```
data(admixed)
str(admixed)
```

---

gen\_structured\_model     *Simulation Scenario from Bhatnagar et al. (2018+) ggmix paper*

---

## Description

Function that generates data of the different simulation studies presented in the accompanying paper. This function requires the popkin and bnpsd package to be installed.

## Usage

```
gen_structured_model(
  n,
  p_design,
  p_kinship,
  k,
  s,
  Fst,
  b0,
  nPC = 10,
  eta,
  sigma2,
  geography = c("ind", "1d", "circ"),
  percent_causal,
  percent_overlap,
  train_tune_test = c(0.6, 0.2, 0.2)
)
```

## Arguments

n	number of observations to simulate
p_design	number of variables in X_test, i.e., the design matrix
p_kinship	number of variable in X_kinship, i.e., matrix used to calculate kinship
k	number of intermediate subpopulations.
s	the desired bias coefficient, which specifies sigma indirectly. Required if sigma is missing
Fst	The desired final FST of the admixed individuals. Required if sigma is missing
b0	the true intercept parameter
nPC	number of principal components to include in the design matrix used for regression adjustment for population structure via principal components. This matrix is used as the input in a standard lasso regression routine, where there are no random effects.
eta	the true eta parameter, which has to be $0 < \eta < 1$
sigma2	the true sigma2 parameter

geography	the type of geography for simulation the kinship matrix. "ind" is independent populations where every individuals is actually unadmixed, "1d" is a 1D geography and "circ" is circular geography. Default: "ind". See the functions in the bnpsd for details on how this data is actually generated.
percent_causal	percentage of p_design that is causal. must be $0 \leq percent_{causal} \leq 1$ . The true regression coefficients are generated from a standard normal distribution.
percent_overlap	this represents the percentage of causal SNPs that will also be included in the calculation of the kinship matrix
train_tune_test	the proportion of sample size used for training tuning parameter selection and testing. default is 60/20/20 split

## Details

The kinship is estimated using the popkin function from the popkin package. This function will multiple that kinship matrix by 2 to give the expected covariance matrix which is subsequently used in the linear mixed models

## Value

A list with the following elements

- ytrain** simulated response vector for training set
- ytune** simulated response vector for tuning parameter selection set
- ytest** simulated response vector for test set
- xtrain** simulated design matrix for training set
- xtune** simulated design matrix for tuning parameter selection set
- xtest** simulated design matrix for testing set
- xtrain\_lasso** simulated design matrix for training set for lasso model. This is the same as xtrain, but also includes the nPC principal components
- xtune\_lasso** simulated design matrix for tuning parameter selection set for lasso model. This is the same as xtune, but also includes the nPC principal components
- xtest** simulated design matrix for testing set for lasso model. This is the same as xtest, but also includes the nPC principal components
- causal** character vector of the names of the causal SNPs
- beta** the vector of true regression coefficients
- kin\_train** 2 times the estimated kinship for the training set individuals
- kin\_tune\_train** The covariance matrix between the tuning set and the training set individuals
- kin\_test\_train** The covariance matrix between the test set and training set individuals
- Xkinship** the matrix of SNPs used to estimate the kinship matrix
- not\_causal** character vector of the non-causal SNPs
- PC** the principal components for population structure adjustment

**See Also**

[admix\\_prop\\_1d\\_linear](#)

**Examples**

```
admixed <- gen_structured_model(n = 100,
                               p_design = 50,
                               p_kinship = 5e2,
                               geography = "1d",
                               percent_causal = 0.10,
                               percent_overlap = "100",
                               k = 5, s = 0.5, Fst = 0.1,
                               b0 = 0, nPC = 10,
                               eta = 0.1, sigma2 = 1,
                               train_tune_test = c(0.8, 0.1, 0.1))

names(admixed)
```

---

ggmix

---

*Fit Linear Mixed Model with Lasso or Group Lasso Regularization*


---

**Description**

Main function to fit the linear mixed model with lasso or group lasso penalty for a sequence of tuning parameters. This is a penalized regression method that accounts for population structure using either the kinship matrix or the factored realized relationship matrix

**Usage**

```
ggmix(
  x,
  y,
  U,
  D,
  kinship,
  K,
  n_nonzero_eigenvalues,
  n_zero_eigenvalues,
  estimation = c("full"),
  penalty = c("lasso"),
  group,
  penalty.factor = rep(1, p_design),
  lambda = NULL,
  lambda_min_ratio = ifelse(n_design < p_design, 0.01, 1e-04),
  nlambdas = 100,
  eta_init = 0.5,
  maxit = 100,
  fdev = 1e-20,
  standardize = FALSE,
```

```

alpha = 1,
thresh_glmnet = 1e-08,
epsilon = 1e-04,
dfmax = p_design + 2,
verbose = 0
)

```

## Arguments

x	input matrix, of dimension n x p; where n is the number of observations and p are the number of predictors.
y	response variable. must be a quantitative variable
U	left singular vectors corresponding to the non-zero eigenvalues provided in the D argument.
D	non-zero eigenvalues. This option is provided to the user should they decide or need to calculate the eigen decomposition of the kinship matrix or the singular value decomposition of the matrix of SNPs used to calculate the kinship outside of this function. This may occur, if for example, it is easier (e.g. because of memory issues, it's easier to calculate in plink). This should correspond to the non-zero eigenvalues only. Note that if you are doing an svd on the matrix of SNPs used to calculate the kinship matrix, then you must provide the square of the singular values so that they correspond to the eigenvalues of the kinship matrix. If you want to use the low rank estimation algorithm, you must provide the truncated eigenvalues and eigenvectors to the D and U arguments, respectively. If you want ggmix to truncate the eigenvectors and eigenvalues for low rank estimation, then provide either K or kinship instead and specify n_nonzero_eigenvalues.
kinship	positive definite kinship matrix
K	the matrix of SNPs used to determine the kinship matrix
n_nonzero_eigenvalues	the number of nonzero eigenvalues. This argument is only used when estimation="low" and either kinship or K is provided. This argument will limit the function to finding the n_nonzero_eigenvalues largest eigenvalues. If U and D have been provided, then n_nonzero_eigenvalues defaults to the length of D.
n_zero_eigenvalues	Currently not being used. Represents the number of zero eigenvalues. This argument must be specified when U and D are specified and estimation="low". This is required for low rank estimation because the number of zero eigenvalues and their corresponding eigenvalues appears in the likelihood. In general this would be the rank of the matrix used to calculate the eigen or singular value decomposition. When kinship is provided and estimation="low" the default value will be nrow(kinship) - n_nonzero_eigenvalues. When K is provided and estimation="low", the default value is rank(K) - n_nonzero_eigenvalues
estimation	type of estimation. Currently only type="full" has been implemented.
penalty	type of regularization penalty. Currently, only penalty="lasso" has been implemented.

group	a vector of consecutive integers describing the grouping of the coefficients. Currently not implemented, but will be used when penalty="gglasso" is implemented.
penalty.factor	Separate penalty factors can be applied to each coefficient. This is a number that multiplies lambda to allow differential shrinkage. Can be 0 for some variables, which implies no shrinkage, and that variable is always included in the model. Default is 1 for all variables
lambda	A user supplied lambda sequence (this is the tuning parameter). Typical usage is to have the program compute its own lambda sequence based on nlambda and lambda.min.ratio. Supplying a value of lambda overrides this. WARNING: use with care. Do not supply a single value for lambda (for predictions after CV use predict() instead). Supply instead a decreasing sequence of lambda values. glmnet relies on its warm starts for speed, and it's often faster to fit a whole path than compute a single fit.
lambda.min.ratio	Smallest value for lambda, as a fraction of lambda.max, the (data derived) entry value (i.e. the smallest value for which all coefficients are zero). The default depends on the sample size nobs relative to the number of variables nvars. If nobs > nvars, the default is 0.0001, close to zero. If nobs < nvars, the default is 0.01. A very small value of lambda.min.ratio will lead to a saturated fit in the nobs < nvars case.
nlambda	the number of lambda values - default is 100.
eta_init	initial value for the eta parameter, with $0 < \eta < 1$ used in determining lambda.max and starting value for fitting algorithm.
maxit	Maximum number of passes over the data for all lambda values; default is $10^2$ .
fdev	Fractional deviance change threshold. If change in deviance between adjacent lambdas is less than fdev, the solution path stops early. factory default = $1.0e-5$
standardize	Logical flag for x variable standardization, prior to fitting the model sequence. The coefficients are always returned on the original scale. Default is standardize=FALSE. If variables are in the same units already, you might not wish to standardize.
alpha	The elasticnet mixing parameter, with $0 \leq \alpha \leq 1$ . alpha=1 is the lasso penalty, and alpha=0 the ridge penalty.
thresh_glmnet	Convergence threshold for coordinate descent for updating beta parameters. Each inner coordinate-descent loop continues until the maximum change in the objective after any coefficient update is less than thresh times the null deviance. Defaults value is $1E-7$
epsilon	Convergence threshold for block relaxation of the entire parameter vector $\Theta = (\beta, \eta, \sigma^2)$ . The algorithm converges when <div style="text-align: center;"> <math display="block">crossprod(\Theta_{j+1} - \Theta_j) &lt; \epsilon</math> </div> . Defaults value is $1E-4$
dfmax	limit the maximum number of variables in the model. Useful for very large p (the total number of predictors in the design matrix), if a partial path is desired. Default is the number of columns in the design matrix + 2 (for the variance components)



verbose            display progress. Can be either 0,1 or 2. 0 will not display any progress, 2 will display very detailed progress and 1 is somewhere in between. Default: 0.

### Value

list which includes the fitted object, lambda sequence, solution path, variance-covariance parameters, degrees of freedom, and singular values, vectors of kinship matrix

### References

Bhatnagar, Sahir R, Yang, Yi, Lu, Tianyuan, Schurr, Erwin, Loredó-Osti, JC, Forest, Marie, Oualkacha, Karim, and Greenwood, Celia MT. (2020) *Simultaneous SNP selection and adjustment for population structure in high dimensional prediction models* doi: [10.1101/408484](https://doi.org/10.1101/408484)

Friedman, J., Hastie, T. and Tibshirani, R. (2008) *Regularization Paths for Generalized Linear Models via Coordinate Descent*, <https://web.stanford.edu/~hastie/Papers/glmnet.pdf> *Journal of Statistical Software*, Vol. 33(1), 1-22 Feb 2010 <https://www.jstatsoft.org/v33/i01/>

Yang, Y., & Zou, H. (2015). A fast unified algorithm for solving group-lasso penalized learning problems. *Statistics and Computing*, 25(6), 1129-1141. [https://www.math.mcgill.ca/yyang/resources/papers/STCO\\_gglasso.pdf](https://www.math.mcgill.ca/yyang/resources/papers/STCO_gglasso.pdf)

### Examples

```
data(admixed)
fitlmm <- ggmix(x = admixed$xtrain, y = admixed$ytrain,
               kinship = admixed$kin_train,
               estimation = "full")
gicfit <- gic(fitlmm)
coef(gicfit, type = "nonzero")
predict(gicfit, newx = admixed$xtest)[1:5,drop=FALSE]
plot(gicfit)
plot(fitlmm)
```

---

ggmix\_data\_object

---

*Constructor functions for the different ggmix objects*


---

### Description

new\_fullrank\_kinship, new\_fullrank\_K, new\_fullrank\_UD, new\_lowrank\_kinship, new\_lowrank\_K and new\_lowrank\_UD create the ggmix objects from the provided data that are necessary to fit the penalized linear mixed model according to the user's parameters.

### Usage

```
new_fullrank_kinship(x, y, kinship)
```

```
new_fullrank_K(x, y, K)
```

```

new_fullrank_UD(x, y, U, D)

new_lowrank_kinship(x, y, kinship, n_nonzero_eigenvalues, n_zero_eigenvalues)

new_lowrank_K(x, y, K, n_nonzero_eigenvalues, n_zero_eigenvalues)

new_lowrank_UD(x, y, U, D, n_nonzero_eigenvalues, n_zero_eigenvalues)

```

### Arguments

x	input matrix, of dimension $n \times p$ ; where $n$ is the number of observations and $p$ are the number of predictors.
y	response variable. must be a quantitative variable
kinship	positive definite kinship matrix
K	the matrix of SNPs used to determine the kinship matrix
U	left singular vectors corresponding to the non-zero eigenvalues provided in the D argument.
D	non-zero eigenvalues. This option is provided to the user should they decide or need to calculate the eigen decomposition of the kinship matrix or the singular value decomposition of the matrix of SNPs used to calculate the kinship outside of this function. This may occur, if for example, it is easier (e.g. because of memory issues, it's easier to calculate in <code>plink</code> ). This should correspond to the non-zero eigenvalues only. Note that if you are doing an <code>svd</code> on the matrix of SNPs used to calculate the kinship matrix, then you must provide the square of the singular values so that they correspond to the eigenvalues of the kinship matrix. If you want to use the low rank estimation algorithm, you must provide the truncated eigenvalues and eigenvectors to the D and U arguments, respectively. If you want <code>ggmix</code> to truncate the eigenvectors and eigenvalues for low rank estimation, then provide either K or kinship instead and specify <code>n_nonzero_eigenvalues</code> .
<code>n_nonzero_eigenvalues</code>	the number of nonzero eigenvalues. This argument is only used when <code>estimation="low"</code> and either kinship or K is provided. This argument will limit the function to finding the <code>n_nonzero_eigenvalues</code> largest eigenvalues. If U and D have been provided, then <code>n_nonzero_eigenvalues</code> defaults to the length of D.
<code>n_zero_eigenvalues</code>	the number of desired or specified zero eigenvalues. This is only needed when <code>estimation="lowrank"</code> , and is calculated internally by the <code>ggmix</code> function. It is equal to the number of observations minus <code>n_nonzero_eigenvalues</code>

### Value

A `ggmix` object, of the class that corresponds to the estimation method. These objects are lists that contain the data necessary for computation. These functions are not meant to be called directly by the user

**See Also**[ggmix](#)

gic

*Generalised Information Criterion***Description**

Calculates the generalised information criterion for each value of the tuning parameter lambda

**Usage**

```
gic(ggmix_fit, ...)

## Default S3 method:
gic(ggmix_fit, ...)

## S3 method for class 'ggmix_fit'
gic(ggmix_fit, ..., an = log(log(n)) * log(p))
```

**Arguments**

**ggmix\_fit** An object of class ggmix\_fit which is outputted by the [ggmix](#) function

**...** other parameters. currently ignored.

**an** numeric, the penalty per parameter to be used; the default is  $an = \log(\log(n)) * \log(p)$  where n is the number of subjects and p is the number of parameters

**Details**

the generalised information criterion used for gaussian response is given by

$$-2 * \loglikelihood(\hat{\Theta}) + an * df$$

where df is the number of non-zero estimated parameters, including variance components

**Value**

an object with S3 class "ggmix\_gic", "ggmix\_fit", "\*" and "\*\*\*" where "\*" is "lasso" or "gglasso" and "\*\*\*" is fullrank or lowrank. Results are provided for converged values of lambda only.

**ggmix\_fit** the ggmix\_fit object

**lambda** the sequence of converged tuning parameters

**nzero** the number of non-zero estimated coefficients including the 2 variance parameters which are not penalized and therefore always included

**gic** gic value. a numeric vector with length equal to length(lambda)

**lambda.min.name** a character corresponding to the name of the tuning parameter lambda which minimizes the gic

**lambda.min** the value of lambda which minimizes the gic

**References**

Fan Y, Tang CY. Tuning parameter selection in high dimensional penalized likelihood. Journal of the Royal Statistical Society: Series B (Statistical Methodology). 2013 Jun 1;75(3):531-52.

Nishii R. Asymptotic properties of criteria for selection of variables in multiple regression. The Annals of Statistics. 1984;12(2):758-65.

**See Also**

[ggmix](#)

---

gr\_eta\_lasso\_fullrank *Functions related to eta parameter used in optim and kkt checks*

---

**Description**

Used for gradient of eta. Currently being passed to optim in [lmmlasso](#) and used in [kkt\\_check](#)

**Usage**

```
gr_eta_lasso_fullrank(eta, sigma2, beta, eigenvalues, x, y, nt)
```

```
fn_eta_lasso_fullrank(eta, sigma2, beta, eigenvalues, x, y, nt)
```

**Arguments**

eta	current estimate of the eta parameter
sigma2	current estimate of the sigma2 parameter
beta	current estimate of the beta parameter including the intercept. this should be of length p+1, where p is the number of variables.
eigenvalues	non-zero eigenvalues of the kinship matrix, or the square of the singular values of the matrix used to construct the kinship matrix
x	input matrix, of dimension n x p; where n is the number of observations and p are the number of predictors.
y	response variable. must be a quantitative variable
nt	total number of observations

**Value**

returns the value of the function and gradient for the eta variance parameter

**See Also**

[logliklasso](#), [kkt\\_check](#), [lmmlasso](#)

---

karim	<i>Karim's Simulated Data</i>
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---

## Description

A simulated dataset with a kinship matrix

## Usage

```
karim
```

## Format

A list with 6 elements:

**b** vector of length 1000 representing the true regression coefficients. 10 non-zero coefficients, the rest are 0.

**kin1** the true kinship matrix

**s.g** polygenic variance, set to be 1.26

**s.e** error variance, set to be 1

**h.tot** the total trait heritability. Set to be 60 of genotypes of dimension 600 x 1000 SNPs, with approximately 800 common and 200 rare SNPs

## Details

If you simulate data using the scenario provided in the example, then the QTL heritability of y will be 8 of the trait's total heritability), and the trait total heritability is set to be 60

## Examples

```
data(karim)
# Simulate a response using the genotype matrix and the kinship matrix
Phi <- 2 * karim$kin1
intercept <- 1
P <- MASS::mvrnorm(1, rep(0,600), karim$s.g * Phi)
y <- intercept + karim$G %*% karim$b + P + rnorm(600,0,karim$s.e)
```

kkt\_check

*Check of KKT Conditions for Linear Mixed Model***Description**

This function checks the KKT conditions

**Usage**

```
kkt_check(eta, sigma2, beta, eigenvalues, x, y, nt, lambda, tol.kkt = 0.001)
```

```
grr_sigma2(eta, sigma2, beta, eigenvalues, x, y, nt)
```

```
grr_beta0(eta, sigma2, beta, eigenvalues, x, y, nt)
```

**Arguments**

eta	current estimate of the eta parameter
sigma2	current estimate of the sigma2 parameter
beta	current estimate of the beta parameter including the intercept. this should be of length $p+1$ , where $p$ is the number of variables.
eigenvalues	non-zero eigenvalues of the kinship matrix, or the square of the singular values of the matrix used to construct the kinship matrix
x	rotated x. Should be $U^T X$ , where $U$ is the matrix of eigenvectors and $X$ contains the first column of ones for the intercept. $x$ should be a matrix of dimension $n \times (p+1)$ . These are outputted by the constructor functions. See <a href="#">ggmix_data_object</a> for details
y	rotated y. Should be $U^T Y$ , where $U$ is the matrix of eigenvectors and $Y$ is the response.
nt	total number of observations
lambda	A user supplied lambda sequence (this is the tuning parameter). Typical usage is to have the program compute its own lambda sequence based on <code>nlambdas</code> and <code>lambda.min.ratio</code> . Supplying a value of lambda overrides this. <b>WARNING:</b> use with care. Do not supply a single value for lambda (for predictions after CV use <code>predict()</code> instead). Supply instead a decreasing sequence of lambda values. <code>glmnet</code> relies on its warm starts for speed, and its often faster to fit a whole path than compute a single fit.
tol.kkt	Tolerance for determining if an entry of the subgradient is zero

**Value**

returns the values of the gradient for each of the parameters

**Note**

`grr_sigma2` and `grr_beta0` are functions for the gradient of `sigma2` and `beta0`, respectively

---

lambdasso	<i>Estimation of Lambda Sequence for Linear Mixed Model with Lasso Penalty</i>
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---

## Description

lambdasso estimates a decreasing sequence of tuning parameters

## Usage

```
lambdasso(ggmix_object, ...)

## Default S3 method:
lambdasso(ggmix_object, ...)

## S3 method for class 'fullrank'
lambdasso(
  ggmix_object,
  ...,
  penalty.factor,
  lambda_min_ratio,
  epsilon = 1e-14,
  tol.kkt = 1e-09,
  eta_init = 0.5,
  nlambdas = 100,
  scale_x = FALSE,
  center_y = FALSE
)
```

## Arguments

ggmix_object	A ggmix_object object of class lowrank or fullrank
...	Extra parameters. Currently ignored.
penalty.factor	Separate penalty factors can be applied to each coefficient. This is a number that multiplies lambda to allow differential shrinkage. Can be 0 for some variables, which implies no shrinkage, and that variable is always included in the model. Default is 1 for all variables
lambda_min_ratio	Smallest value for lambda, as a fraction of lambda.max, the (data derived) entry value (i.e. the smallest value for which all coefficients are zero). The default depends on the sample size nobs relative to the number of variables nvars. If nobs > nvars, the default is 0.0001, close to zero. If nobs < nvars, the default is 0.01. A very small value of lambda.min.ratio will lead to a saturated fit in the nobs < nvars case.

epsilon	Convergence threshold for block relaxation of the entire parameter vector $\Theta = (\beta, \eta, \sigma^2)$ . The algorithm converges when $\text{crossprod}(\Theta_{j+1} - \Theta_j) < \epsilon$ . Defaults value is 1E-4
tol.kkt	KKT tolerance. Currently ignored
eta_init	initial value for the eta parameter, with $0 < \eta < 1$ used in determining lambda.max and starting value for fitting algorithm.
nlambda	the number of lambda values - default is 100.
scale_x	should the columns of x be scaled - default is FALSE
center_y	should y be mean centered - default is FALSE.

**Value**

A decreasing sequence of tuning parameters

**Note**

This function isn't meant to be called directly by the user.

**See Also**

[ggmix](#)

---

lmmlasso

*Estimation of Linear Mixed Model with Lasso Penalty*


---

**Description**

lmmlasso estimates the linear mixed model with lasso penalty

**Usage**

```
lmmlasso(ggmix_object, ...)

## Default S3 method:
lmmlasso(ggmix_object, ...)

## S3 method for class 'fullrank'
lmmlasso(
  ggmix_object,
  ...,
  penalty.factor,
  lambda,
  lambda_min_ratio,
```



```

    nlambda,
    n_design,
    p_design,
    eta_init,
    maxit,
    fdev,
    standardize,
    alpha,
    thresh_glmnet,
    epsilon,
    dfmax,
    verbose
)

```

## Arguments

<code>ggmix_object</code>	A <code>ggmix_object</code> object of class <code>lowrank</code> or <code>fullrank</code>
<code>...</code>	Extra parameters. Currently ignored.
<code>penalty.factor</code>	Separate penalty factors can be applied to each coefficient. This is a number that multiplies <code>lambda</code> to allow differential shrinkage. Can be 0 for some variables, which implies no shrinkage, and that variable is always included in the model. Default is 1 for all variables
<code>lambda</code>	A user supplied <code>lambda</code> sequence (this is the tuning parameter). Typical usage is to have the program compute its own <code>lambda</code> sequence based on <code>nlambda</code> and <code>lambda.min.ratio</code> . Supplying a value of <code>lambda</code> overrides this. <b>WARNING:</b> use with care. Do not supply a single value for <code>lambda</code> (for predictions after CV use <code>predict()</code> instead). Supply instead a decreasing sequence of <code>lambda</code> values. <code>glmnet</code> relies on its warm starts for speed, and its often faster to fit a whole path than compute a single fit.
<code>lambda.min.ratio</code>	Smallest value for <code>lambda</code> , as a fraction of <code>lambda.max</code> , the (data derived) entry value (i.e. the smallest value for which all coefficients are zero). The default depends on the sample size <code>nobs</code> relative to the number of variables <code>nvars</code> . If <code>nobs &gt; nvars</code> , the default is 0.0001, close to zero. If <code>nobs &lt; nvars</code> , the default is 0.01. A very small value of <code>lambda.min.ratio</code> will lead to a saturated fit in the <code>nobs &lt; nvars</code> case.
<code>nlambda</code>	the number of <code>lambda</code> values - default is 100.
<code>n_design</code>	total number of observations
<code>p_design</code>	number of variables in the design matrix, excluding the intercept column
<code>eta_init</code>	initial value for the <code>eta</code> parameter, with $0 < \eta < 1$ used in determining <code>lambda.max</code> and starting value for fitting algorithm.
<code>maxit</code>	Maximum number of passes over the data for all <code>lambda</code> values; default is $10^2$ .
<code>fdev</code>	Fractional deviance change threshold. If change in deviance between adjacent <code>lambdas</code> is less than <code>fdev</code> , the solution path stops early. factory default = $1.0e-5$

standardize	Logical flag for x variable standardization, prior to fitting the model sequence. The coefficients are always returned on the original scale. Default is standardize=FALSE. If variables are in the same units already, you might not wish to standardize.
alpha	The elasticnet mixing parameter, with $0 \leq \alpha \leq 1$ . alpha=1 is the lasso penalty, and alpha=0 the ridge penalty.
thresh_glmnet	Convergence threshold for coordinate descent for updating beta parameters. Each inner coordinate-descent loop continues until the maximum change in the objective after any coefficient update is less than thresh times the null deviance. Defaults value is 1E-7
epsilon	Convergence threshold for block relaxation of the entire parameter vector $\Theta = (\beta, \eta, \sigma^2)$ . The algorithm converges when <div style="text-align: center;"> <math display="block">crossprod(\Theta_{j+1} - \Theta_j) &lt; \epsilon</math> </div> . Defaults value is 1E-4
dfmax	limit the maximum number of variables in the model. Useful for very large p (the total number of predictors in the design matrix), if a partial path is desired. Default is the number of columns in the design matrix + 2 (for the variance components)
verbose	display progress. Can be either 0,1 or 2. 0 will not display any progress, 2 will display very detailed progress and 1 is somewhere in between. Default: 0.

**Value**

A object of class ggmix

**See Also**

[ggmix](#)

---

logliklasso	<i>Estimation of Log-likelihood for Linear Mixed Model with Lasso Penalty</i>
-------------	---

---

**Description**

sigma2lasso estimates the value of the sigma2 for the linear mixed model with lasso penalty

**Usage**

```
logliklasso(ggmix_object, ...)

## Default S3 method:
logliklasso(ggmix_object, ...)

## S3 method for class 'fullrank'
logliklasso(ggmix_object, ..., eta, sigma2, beta, nt, x, y)
```

**Arguments**

ggmix_object	A ggmix_object object of class lowrank or fullrank
...	Extra parameters. Currently ignored.
eta	current estimate of the eta parameter
sigma2	current estimate of the sigma2 parameter
beta	current estimate of the beta parameter including the intercept. this should be of length $p+1$ , where $p$ is the number of variables.
nt	total number of observations
x	input matrix, of dimension $n \times p$ ; where $n$ is the number of observations and $p$ are the number of predictors.
y	response variable. must be a quantitative variable

**Value**

A decreasing sequence of tuning parameters

**Note**

This function isn't meant to be called directly by the user.

**See Also**

[ggmix](#)  
[ggmix\\_data\\_object](#)

---

plot.ggmix_fit	<i>Plot Method for ggmix_fit object</i>
----------------	---

---

**Description**

Produces a coefficient profile plot of the coefficient paths for a fitted ggmix\_fit object.

**Usage**

```
## S3 method for class 'ggmix_fit'
plot(x, ..., xvar = c("norm", "lambda", "dev"), label = FALSE, sign.lambda = 1)

plotCoef(
  beta,
  norm,
  lambda,
  df,
  dev,
  label = FALSE,
  xvar = c("norm", "lambda", "dev"),
```

```

    xlab = iname,
    ylab = "Coefficients",
    ...
)

```

### Arguments

x	a ggmix_fit object
...	other graphical parameters passed to plot
xvar	What is on the X-axis. "norm" plots against the L1-norm of the coefficients, "lambda" against the log-lambda sequence, and "dev" against the percent deviance explained.
label	If TRUE, label the curves with variable sequence numbers.
sign.lambda	Either plot against log(lambda) (default) or its negative if sign.lambda=-1
beta	fixed effects estimates
norm	l1 norm of fixed effect estimates. if missing, (default) this function will calculate it
lambda	sequence of tuning parameters
df	number of non-zero fixed + random effects
dev	percent deviance
xlab	x-axis label
ylab	y-axis label

### Details

A coefficient profile plot is produced

### Value

A plot is produced and nothing is returned

---

plot.ggmix\_gic

*Plot the Generalised Information Criteria curve produced by gic*

---

### Description

Plots the Generalised Information Criteria curve, as a function of the lambda values used

**Usage**

```
## S3 method for class 'ggmix_gic'
plot(
  x,
  ...,
  sign.lambda = 1,
  type = c("gic", "QQranef", "QQresid", "predicted", "Tukey-Anscombe"),
  s = "lambda.min",
  newy,
  newx
)

plotGIC(x, sign.lambda, lambda.min, ...)
```

**Arguments**

x	fitted linear mixed model object of class ggmix_gic from the <a href="#">gic</a> function
...	Other graphical parameters to plot
sign.lambda	Either plot against log(lambda) (default) or its negative if sign.lambda=-1
type	gic returns a plot of the GIC vs. log(lambda). QQranef return a qqplot of the random effects. QQresid returns a qqplot of the residuals which is $y - X\beta - b_i$ where b_i is the subject specific random effect. predicted returns a plot of the predicted response ( $X\beta + b_i$ ) vs. the observed response, where b_i is the subject specific random effect. Tukey-Anscombe returns a plot of the residuals vs. fitted values ( $X\beta$ )
s	Value of the penalty parameter lambda at which predictions are required. Default is the value s="lambda.min". If s is numeric, it is taken as the value of lambda to be used. Must be a single value of the penalty parameter lambda at which coefficients will be extracted via the coef method for objects of class ggmix_gic. If more than one is supplied, only the first one will be used.
newy	the response variable that was provided to ggmix. this is only required for type="QQresid", type="Tukey-Anscombe" and type="predicted"
newx	matrix of values for x at which predictions are to be made. Do not include the intercept. this is only required for type="QQresid", type="Tukey-Anscombe" and type="predicted"
lambda.min	the value of lambda which minimizes the gic

**Details**

A plot is produced, and nothing is returned.

**Value**

plot depends on the type selected

**See Also**

[gic](#)

**Examples**

```
data("admixed")
fit <- ggmix(x = admixed$xtrain,
            y = admixed$ytrain,
            kinship = admixed$kin_train)
hdbic <- gic(fit)

# plot solution path
plot(fit)

# plot HDBIC curve as a function of lambda
plot(hdbic)
```

---

predict.ggmix_fit	<i>Make predictions from a ggmix_fit object</i>
-------------------	---

---

**Description**

Similar to other predict methods, this functions predicts fitted values, coefficients and more from a fitted ggmix\_fit object.

**Usage**

```
## S3 method for class 'ggmix_fit'
predict(
  object,
  newx,
  s = NULL,
  type = c("link", "response", "coefficients", "all", "nonzero", "individual"),
  covariance,
  ...
)

## S3 method for class 'ggmix_fit'
coef(object, s = NULL, type, ...)
```

**Arguments**

object	Fitted ggmix_fit model object from the <a href="#">ggmix</a> function
newx	matrix of values for x at which predictions are to be made. Do not include the intercept. Must be a matrix. This argument is not used for type = c("coefficients", "nonzero", "all"). This matrix must have the same number of columns originally supplied to the <a href="#">ggmix</a> fitting function.
s	Value(s) of the penalty parameter lambda at which predictions are required. Default is the entire sequence used to create the model.

type	Type of prediction required. Type "link" gives the fitted values $X\beta$ . Type "response" is equivalent to type "link". Type "coefficients" computes the coefficients at the requested values for s and returns the regression coefficients only, including the intercept. Type "all" returns both the regression coefficients and variance components at the requested value of s. Type "nonzero" returns a 1 column matrix of the the nonzero fixed effects, as well as variance components for each value of s. If more than one s is provided, then "nonzero" will return a list of 1 column matrices. Default: "link"
covariance	covariance between test and training individuals. if there are q testing individuals and N-q training individuals, then this covariance matrix is q x (N-q)
...	additional arguments to pass to predict function

### Details

s is the new vector at which predictions are requested. If s is not in the lambda sequence used for fitting the model, the predict function will use linear interpolation to make predictions. The new values are interpolated using a fraction of predicted values from both left and right lambda indices. `coef(...)` is equivalent to `predict(ggmix_fit, type="coefficients", ...)`. To get individual level predictions at each value of lambda, you must provide the lambda sequence to the s argument. You can pass either a `ggmix_fit` or `ggmix_gic` object. See examples for more details.

### Value

The object returned depends on type.

### Examples

```
data("admixed")
fitlmm <- ggmix(x = admixed$xtrain, y = admixed$ytrain,
               kinship = admixed$kin_train,
               estimation = "full")
bicGGMIX <- gic(fitlmm,
               an = log(length(admixed$ytrain)))
plot(bicGGMIX)
coef(bicGGMIX, s = "lambda.min")
yhat_test <- predict(bicGGMIX, s="lambda.min",
                    newx = admixed$xtest, type = "individual",
                    covariance = admixed$kin_test_train)
cor(yhat_test, admixed$ytest)
yhat_test_population <- predict(bicGGMIX, s="lambda.min",
                              newx = admixed$xtest,
                              type = "response")
```

## Description

This function makes predictions from a `ggmix_gic` object, using the stored `"ggmix_fit"` object, and the optimal value chosen for `lambda` using the `gic`.

## Usage

```
## S3 method for class 'ggmix_gic'
predict(object, newx, s = c("lambda.min"), ...)
```

```
## S3 method for class 'ggmix_gic'
coef(object, s = c("lambda.min"), type, ...)
```

## Arguments

<code>object</code>	fitted <code>ggmix_gic</code> object
<code>newx</code>	matrix of values for <code>x</code> at which predictions are to be made. Do not include the intercept. Must be a matrix. This argument is not used for <code>type = c("coefficients", "nonzero", "all")</code> . This matrix must have the same number of columns originally supplied to the <code>ggmix</code> fitting function.
<code>s</code>	Value(s) of the penalty parameter <code>lambda</code> at which predictions are required. Default is the value <code>s="lambda.min"</code> can be used. If <code>s</code> is numeric, it is taken as the value(s) of <code>lambda</code> to be used.
<code>...</code>	other arguments passed to <code>predict.ggmix_fit</code>
<code>type</code>	Type of prediction required. Type <code>"link"</code> gives the fitted values $X\beta$ . Type <code>"response"</code> is equivalent to type <code>"link"</code> . Type <code>"coefficients"</code> computes the coefficients at the requested values for <code>s</code> and returns the regression coefficients only, including the intercept. Type <code>"all"</code> returns both the regression coefficients and variance components at the requested value of <code>s</code> . Type <code>"nonzero"</code> returns a 1 column matrix of the the nonzero fixed effects, as well as variance components for each value of <code>s</code> . If more than one <code>s</code> is provided, then <code>"nonzero"</code> will return a list of 1 column matrices. Default: <code>"link"</code>

## Details

This function makes it easier to use the results of `gic` chosen model to make a prediction.

## Value

The object returned depends the `...` argument which is passed on to the `predict` method for `ggmix_fit` objects.

## See Also

[predict.ggmix\\_fit](#)



---

print.ggmix_fit	<i>Print Method for Objects of Class ggmix_fit</i>
-----------------	--

---

## Description

print method for objects of class ggmix\_fit

## Usage

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

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

## Arguments

x	an object of class objects of class ggmix_fit
...	other arguments passed to print
digits	significant digits in printout. Default: max(3,getOption("digits") - 3)

## Value

The call that produced the object x is printed, followed by a three-column matrix with columns Df, %Dev, and and Lambda. The Df columns correspond to the number of nonzero coefficients including variance components. %dev is the percent deviance explained (relative to the null deviance). Lambda is the sequence of converged tuning parameters.

## See Also

[ggmix](#)

---

ranef	<i>Extract Random Effects</i>
-------	-------------------------------

---

## Description

Generic function for extracting the random effects. This is the same generic (and same name) defined in the nlme and lme4 package.

**Usage**

```

ranef(object, ...)

random.effects(object, ...)

## Default S3 method:
random.effects(object, ...)

## Default S3 method:
ranef(object, ...)

## S3 method for class 'ggmix_gic'
ranef(object, s = "lambda.min", ...)

```

**Arguments**

<code>object</code>	any fitted model object from which random effects estimates can be extracted. Currently supports "ggmix_gic" objects outputted by the <a href="#">gic</a> function
<code>...</code>	other parameters. currently ignored
<code>s</code>	Value(s) of the penalty parameter lambda at which predictions are required. Default is the value <code>s="lambda.min"</code> can be used. If <code>s</code> is numeric, it is taken as the value(s) of lambda to be used.

**Details**

For objects of class "ggmix\_gic", this function returns the subject-specific random effect value for the model which minimizes the GIC using the maximum a posteriori principle

**Value**

a numeric vector of length equal to the number of observations of subject-specific random effects

**See Also**

[gic](#)

**Examples**

```

data("admixed")
fit <- ggmix(x = admixed$xtrain, y = admixed$ytrain,
            kinship = admixed$kin_train)
gicfit <- gic(fit)
# random effect at selected value of lambda
plot(ggmix::ranef(gicfit))
# random effects at specific values of lambda
head(ggmix::ranef(gicfit, s = c(0.1,0.2)))

```

**Description**

sigma2lasso estimates the value of the sigma<sup>2</sup> for the linear mixed model with lasso penalty

**Usage**

```
sigma2lasso(ggmix_object, ...)  
  
## Default S3 method:  
sigma2lasso(ggmix_object, ...)  
  
## S3 method for class 'fullrank'  
sigma2lasso(ggmix_object, ..., n, beta, eta)
```

**Arguments**

ggmix_object	A ggmix_object object of class lowrank or fullrank
...	Extra parameters. Currently ignored.
n	number of observations
beta	current estimate of the beta parameter including the intercept. this should be of length p+1, where p is the number of variables.
eta	current estimate of the eta parameter

**Value**

A decreasing sequence of tuning parameters

**Note**

There is a closed form solution for sigma<sup>2</sup>, given beta and eta. This function isn't meant to be called directly by the user.

**See Also**

[ggmix](#)

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