

# Package ‘vennLasso’

October 12, 2022

**Type** Package

**Title** Variable Selection for Heterogeneous Populations

**Version** 0.1.6

**Description** Provides variable selection and estimation routines for models with main effects stratified on multiple binary factors. The 'vennLasso' package is an implementation of the method introduced in Huling, et al. (2017) <[doi:10.1111/biom.12769](https://doi.org/10.1111/biom.12769)>.

**URL** <https://github.com/jaredhuling/vennLasso>

**BugReports** <https://github.com/jaredhuling/vennLasso/issues>

**License** GPL (>= 2)

**Encoding** UTF-8

**LazyData** TRUE

**Depends** R (>= 3.2.0)

**Imports** Rcpp (>= 0.11.0), foreach, survival, MASS, Matrix, VennDiagram, visNetwork, igraph, methods

**LinkingTo** Rcpp, RcppEigen, RcppNumerical

**RoxygenNote** 7.1.0

**Suggests** knitr, rmarkdown

**VignetteBuilder** knitr

**NeedsCompilation** yes

**Author** Jared Huling [aut, cre] (<<https://orcid.org/0000-0003-0670-4845>>),  
Muxuan Liang [ctb],  
Yixuan Qiu [cph],  
Gael Guennebaud [cph],  
Ray Gardner [cph],  
Jitse Niesen [cph]

**Maintainer** Jared Huling <jaredhuling@gmail.com>

**Repository** CRAN

**Date/Publication** 2020-06-03 08:50:02 UTC

## R topics documented:

cv.vennLasso	2
estimate.hier.sparsity.param	5
genHierSparseBeta	6
genHierSparseData	7
logLik.vennLasso	10
oglasso	11
plot.cv.vennLasso	14
plotCoefs	15
plotSelections	16
plotVenn	17
predict.cv.vennLasso	18
predict.vennLasso	19
vennLasso	20

<b>Index</b>	<b>25</b>
--------------	-----------

---

cv.vennLasso	<i>Cross Validation for the vennLasso</i>
--------------	---

---

### Description

Cross Validation for the vennLasso

### Usage

```
cv.vennLasso(
  x,
  y,
  groups,
  lambda = NULL,
  compute.se = FALSE,
  conf.int = NULL,
  type.measure = c("mse", "deviance", "class", "auc", "mae", "brier"),
  nfolds = 10,
  foldid,
  grouped = TRUE,
  keep = FALSE,
  parallel = FALSE,
  ...
)
```

### Arguments

x	input matrix or SparseMatrix of dimension nobs x nvars. Each row is an observation, each column corresponds to a covariate
y	numeric response vector of length nobs

groups	A list of length equal to the number of groups containing vectors of integers indicating the variable IDs for each group. For example, groups=list(c(1,2), c(2,3), c(3,4,5)) specifies that Group 1 contains variables 1 and 2, Group 2 contains variables 2 and 3, and Group 3 contains variables 3, 4, and 5. Can also be a matrix of 0s and 1s with the number of columns equal to the number of groups and the number of rows equal to the number of variables. A value of 1 in row i and column j indicates that variable i is in group j and 0 indicates that variable i is not in group j.
lambda	A user-specified sequence of lambda values. Left unspecified, the a sequence of lambda values is automatically computed, ranging uniformly on the log scale over the relevant range of lambda values.
compute.se	logical flag. If TRUE, standard errors will be computed, otherwise if FALSE they will not
conf.int	value between 0 and 1 indicating the level of the confidence intervals to be computed. For example if conf.int = 0.95, 95 percent confidence intervals will be computed.
type.measure	One of c("mse", "deviance", "class", "auc", "mae", "brier") indicating measure to evaluate for cross-validation. The default is type.measure = "deviance", which uses squared-error for gaussian models (a.k.a type.measure = "mse" there), deviance for logistic regression. type.measure = "class" applies to binomial only. type.measure = "auc" is for two-class logistic regression only. type.measure = "mse" or type.measure = "mae" (mean absolute error) can be used by all models; they measure the deviation from the fitted mean to the response. type.measure = "brier" is for models with family = "coxph" and will compute the Brier score.
nfolds	number of folds for cross-validation. default is 10. 3 is smallest value allowed.
foldid	an optional vector of values between 1 and nfold specifying which fold each observation belongs to.
grouped	Like in <b>glmnet</b> , this is an experimental argument, with default TRUE, and can be ignored by most users. For all models, this refers to computing nfolds separate statistics, and then using their mean and estimated standard error to describe the CV curve. If grouped = FALSE, an error matrix is built up at the observation level from the predictions from the nfold fits, and then summarized (does not apply to type.measure = "auc").
keep	If keep = TRUE, a prevalidated list of array is returned containing fitted values for each observation and each value of lambda for each model. This means these fits are computed with this observation and the rest of its fold omitted. The folid vector is also returned. Default is keep = FALSE
parallel	If TRUE, use parallel foreach to fit each fold. Must register parallel before hand, such as <b>doMC</b> .
...	parameters to be passed to vennLasso

**Value**

An object with S3 class "cv.vennLasso"

**Examples**

```

library(Matrix)

set.seed(123)
n.obs <- 150
n.vars <- 25

true.beta.mat <- array(NA, dim = c(3, n.vars))
true.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
true.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
true.beta.mat[3,] <- c(0, 0, 1, 1, -1, rep(0, n.vars - 5))
rownames(true.beta.mat) <- c("1,0", "1,1", "0,1")
true.beta <- as.vector(t(true.beta.mat))

x.sub1 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub2 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub3 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)

x <- as.matrix(rbind(x.sub1, x.sub2, x.sub3))

conditions <- as.matrix(cbind(c(rep(1, 2 * n.obs), rep(0, n.obs)),
                             c(rep(0, n.obs), rep(1, 2 * n.obs))))

y <- rnorm(n.obs * 3, sd = 3) + drop(as.matrix(bdiag(x.sub1, x.sub2, x.sub3)) %*% true.beta)

fit <- cv.vennLasso(x = x, y = y, groups = conditions, nfolds = 3)

fitted.coef <- predict(fit$vennLasso.fit, type = "coefficients", s = fit$lambda.min)
(true.coef <- true.beta.mat[match(dimnames(fit$vennLasso.fit$beta)[[1]],
                                rownames(true.beta.mat)),])

round(fitted.coef, 2)

## effects smaller for logistic regression
## Not run:
true.beta.mat <- true.beta.mat / 2
true.beta <- true.beta / 2
# logistic regression example#'
y <- rbinom(n.obs * 3, 1,
           prob = 1 / (1 + exp(-drop(as.matrix(bdiag(x.sub1, x.sub2, x.sub3)) %*% true.beta))))

bfit <- cv.vennLasso(x = x, y = y, groups = conditions, family = "binomial",
                   nfolds = 3)

fitted.coef <- predict(bfit$vennLasso.fit, type = "coefficients", s = bfit$lambda.min)
(true.coef <- true.beta.mat[match(dimnames(bfit$vennLasso.fit$beta)[[1]],
                                rownames(true.beta.mat)),])

round(fitted.coef, 2)

## End(Not run)

```

---

```
estimate.hier.sparsity.param
    function to estimate the hierarchical sparsity parameter for a desired
    level of sparsity for simulated hierarchical coefficients
```

---

### Description

function to estimate the hierarchical sparsity parameter for a desired level of sparsity for simulated hierarchical coefficients

### Usage

```
estimate.hier.sparsity.param(
  ncats,
  nvars,
  avg.hier.zeros = 0.3,
  nsims = 150,
  effect.size.max = 0.5,
  misspecification.prop = 0
)
```

### Arguments

<code>ncats</code>	number of categories to stratify on
<code>nvars</code>	number of variables
<code>avg.hier.zeros</code>	desired percent of zero variables among the variables with hierarchical zero patterns.
<code>nsims</code>	number of simulations to estimate the average sparsity. A larger number will be more accurate but take much longer.
<code>effect.size.max</code>	maximum magnitude of the true effect sizes
<code>misspecification.prop</code>	proportion of variables with hierarchical missingness misspecified

### Examples

```
set.seed(123)

# estimate hier.sparsity.param for 0.15 total proportion of nonzero variables
# among vars with hierarchical zero patterns
## Not run:
hsp <- estimate.hier.sparsity.param(ncats = 3, nvars = 25, avg.hier.zeros = 0.15, nsims = 100)

## End(Not run)
# the above results in the following value
hsp <- 0.6341772
```

```

# check that this does indeed achieve the desired level of sparsity
mean(replicate(100, mean(genHierSparseBeta(ncats = 3,
                                           nvars = 25, hier.sparsity.param = hsp) != 0) ))

sparseBeta <- genHierSparseBeta(ncats = 3, nvars = 25, hier.sparsity.param = hsp)

## Not run:
hsp2 <- estimate.hier.sparsity.param(ncats = 2, nvars = 100,
                                     avg.hier.zeros = 0.30, nsims = 50) # 0.5778425
hsp3 <- estimate.hier.sparsity.param(ncats = 3, nvars = 100,
                                     avg.hier.zeros = 0.30, nsims = 50) # 0.4336312
hsp4 <- estimate.hier.sparsity.param(ncats = 4, nvars = 100,
                                     avg.hier.zeros = 0.30, nsims = 50) # 0.2670061
hsp5 <- estimate.hier.sparsity.param(ncats = 5, nvars = 100,
                                     avg.hier.zeros = 0.30, nsims = 50) # 0.146682

## End(Not run)
# 0.07551241 for hsp6

```

---

genHierSparseBeta      *function to generate coefficient matrix with hierarchical sparsity*

---

## Description

function to generate coefficient matrix with hierarchical sparsity

## Usage

```

genHierSparseBeta(
  ncats,
  nvars,
  hier.sparsity.param = 0.5,
  avg.hier.zeros = NULL,
  effect.size.max = 0.5,
  misspecification.prop = 0
)

```

## Arguments

ncats	number of categories to stratify on
nvars	number of variables
hier.sparsity.param	parameter between 0 and 1 which determines how much hierarchical sparsity there is. To achieve a desired total level of sparsity among the variables with hierarchical sparsity, this parameter can be estimated using the function 'estimate.hier.sparsity.param'

`avg.hier.zeros` desired percent of zero variables among the variables with hierarchical zero patterns. If this is specified, it will override the given `hier.sparsity.param` value and estimate it. This takes a while  
`effect.size.max` maximum magnitude of the true effect sizes  
`misspecification.prop` proportion of variables with hierarchical missingness misspecified

## Examples

```

set.seed(123)

# estimate hier.sparsity.param for 0.15 total proportion of nonzero variables
# among vars with hierarchical zero patterns
# NOT RUN: Takes a long time
# hsp <- estimate.hier.sparsity.param(ncats = 3, nvars = 25, avg.hier.zeros = 0.15, nsims = 100)
# the above results in the following value
hsp <- 0.6341772

# check that this does indeed achieve the desired level of sparsity
mean(replicate(100, mean(genHierSparseBeta(ncats = 3,
                                           nvars = 25, hier.sparsity.param = hsp) != 0) ))

sparseBeta <- genHierSparseBeta(ncats = 3, nvars = 25, hier.sparsity.param = hsp)
  
```

---

genHierSparseData      *function to generate data with hierarchical sparsity*

---

## Description

function to generate data with hierarchical sparsity

## Usage

```

genHierSparseData(
  ncats,
  nvars,
  nobs,
  nobs.test = 100,
  hier.sparsity.param = 0.5,
  avg.hier.zeros = NULL,
  prop.zero.vars = 0.5,
  effect.size.max = 0.5,
  misspecification.prop = 0,
  family = c("gaussian", "binomial", "coxph"),
  sd = 1,
  snr = NULL,
  )
  
```

```

    beta = NULL,
    tau = 10,
    covar = 0
  )

```

### Arguments

<code>ncats</code>	number of categories to stratify on
<code>nvars</code>	number of variables
<code>nobs</code>	number of observations per strata to simulate
<code>nobs.test</code>	number of independent test observations per strata to simulate
<code>hier.sparsity.param</code>	parameter between 0 and 1 which determines how much hierarchical sparsity there is. To achieve a desired total level of sparsity among the variables with hierarchical sparsity, this parameter can be estimated using the function <code>'estimate.hier.sparsity.param'</code>
<code>avg.hier.zeros</code>	desired percent of zero variables among the variables with hierarchical zero patterns. If this is specified, it will override the given <code>hier.sparsity.param</code> value and estimate it. This takes a while
<code>prop.zero.vars</code>	proportion of all variables that will be zero across all strata
<code>effect.size.max</code>	maximum magnitude of the true effect sizes
<code>misspecification.prop</code>	proportion of variables with hierarchical missingness misspecified
<code>family</code>	family for the response variable
<code>sd</code>	standard deviation for gaussian simulations
<code>snr</code>	signal-to-noise ratio (only used for <code>family = "gaussian"</code> )
<code>beta</code>	a matrix of true beta values. If given, then no beta will be created and data will be simulated from the given beta
<code>tau</code>	rate parameter for <code>rexp()</code> for generating time-to-event outcomes
<code>covar</code>	scalar, pairwise covariance term for covariates

### Examples

```

set.seed(123)

dat.sim <- genHierSparseData(ncats = 3, nvars = 100, nobs = 200)

# estimate hier.sparsity.param for 0.15 total proportion of nonzero variables
# among vars with hierarchical zero patterns
## Not run:
hsp <- estimate.hier.sparsity.param(ncats = 3, nvars = 50, avg.hier.zeros = 0.15, nsims = 100)

## End(Not run)
# the above results in the following value
hsp <- 0.6270698

```



```
# check that this does indeed achieve the desired level of sparsity
mean(replicate(50, mean(genHierSparseBeta(ncats = 3,
                                       nvars = 50, hier.sparsity.param = hsp) != 0) ))

dat.sim2 <- genHierSparseData(ncats = 3, nvars = 100, nobs = 200, hier.sparsity.param = hsp)

sparseBeta <- genHierSparseBeta(ncats = 3, nvars = 100, hier.sparsity.param = hsp)

## generate data with already generated beta
dat.sim3 <- genHierSparseData(ncats = 3, nvars = 100, nobs = 200, beta = sparseBeta)

## complete example:
## 50% sparsity:
hsp <- 0.2626451

dat.sim <- genHierSparseData(ncats = 3, nvars = 25,
                            nobs = 150, nobs.test = 1000,
                            hier.sparsity.param = hsp,
                            prop.zero.vars = 0.5,
                            effect.size.max = 0.25,
                            family = "gaussian")

x      <- dat.sim$x
x.test <- dat.sim$x.test
y      <- dat.sim$y
y.test <- dat.sim$y.test
grp    <- dat.sim$group.ind
grp.test <- dat.sim$group.ind.test

fit.adapt <- cv.vennLasso(x, y,
                        grp,
                        adaptive.lasso = TRUE,
                        nlambdas      = 25,
                        family        = "gaussian",
                        abs.tol       = 1e-5,
                        rel.tol       = 1e-5,
                        maxit         = 1000,
                        irls.maxit    = 15L,
                        gamma         = 0.2,
                        standardize    = FALSE,
                        intercept     = TRUE,
                        nfolds         = 3,
                        model.matrix  = TRUE)

preds.a <- predict(fit.adapt$vennLasso.fit, x.test, grp.test, s = fit.adapt$lambda.min,
                  type = 'response')
```

---

logLik.vennLasso      *log likelihood function for fitted vennLasso objects*

---

## Description

log likelihood function for fitted vennLasso objects

## Usage

```
## S3 method for class 'vennLasso'
logLik(object, ...)
```

## Arguments

object	fitted "vennLasso" model object.
...	not used

## Examples

```
library(Matrix)

set.seed(123)
n.obs <- 200
n.vars <- 50

true.beta.mat <- array(NA, dim = c(3, n.vars))
true.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
true.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
true.beta.mat[3,] <- c(0, 0, 1, 1, -1, rep(0, n.vars - 5))
rownames(true.beta.mat) <- c("1,0", "1,1", "0,1")
true.beta <- as.vector(t(true.beta.mat))

x.sub1 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub2 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub3 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)

x <- as.matrix(rbind(x.sub1, x.sub2, x.sub3))

conditions <- as.matrix(cbind(c(rep(1, 2 * n.obs), rep(0, n.obs)),
                             c(rep(0, n.obs), rep(1, 2 * n.obs))))

y <- rnorm(n.obs * 3, sd = 3) + drop(as.matrix(bdiag(x.sub1, x.sub2, x.sub3)) %*% true.beta)

fit <- vennLasso(x = x, y = y, groups = conditions)

logLik(fit)
```

---

 oglasso *Overlapping Group Lasso (OGLasso)*


---

**Description**

Overlapping Group Lasso (OGLasso)

**Usage**

```
oglasso(
  x,
  y,
  delta = NULL,
  group,
  fused = NULL,
  family = c("gaussian", "binomial", "coxph"),
  nlambda = 100L,
  lambda = NULL,
  lambda.min.ratio = NULL,
  lambda.fused = 0,
  alpha = NULL,
  group.weights = NULL,
  adaptive.lasso = FALSE,
  adaptive.fused = FALSE,
  penalty.factor = NULL,
  penalty.factor.fused = NULL,
  gamma = 1,
  standardize = TRUE,
  intercept = TRUE,
  compute.se = FALSE,
  rho = NULL,
  dynamic.rho = TRUE,
  maxit = 500L,
  abs.tol = 1e-05,
  rel.tol = 1e-05,
  irls.tol = 1e-05,
  irls.maxit = 100L
)
```

**Arguments**

x	input matrix of dimension nobs by nvars. Each row is an observation, each column corresponds to a covariate
y	numeric response vector of length nobs
delta	vector of length equal to the number of observations with values in 1 and 0, where a 1 indicates the observed time is a death and a 0 indicates the observed time is a censoring event

group	A list of length equal to the number of groups containing vectors of integers indicating the variable IDs for each group. For example, <code>group = list(c(1, 2), c(2, 3), c(3, 4, 5))</code> specifies that Group 1 contains variables 1 and 2, Group 2 contains variables 2 and 3, and Group 3 contains variables 3, 4, and 5. Can also be a matrix of 0s and 1s with the number of columns equal to the number of groups and the number of rows equal to the number of variables. A value of 1 in row <i>i</i> and column <i>j</i> indicates that variable <i>i</i> is in group <i>j</i> and 0 indicates that variable <i>i</i> is not in group <i>j</i> .
fused	matrix specifying generalized lasso penalty formulation. Each column corresponds to each variable and each row corresponds to a new penalty term, ie if row 1 has the first entry of 1 and the second entry of -1, then the penalty term <code>lambda.fused * lbeta_1 - beta_2l</code> will be added. Not available now
family	"gaussian" for least squares problems, "binomial" for binary response
nlambda	The number of lambda values. Default is 100.
lambda	A user-specified sequence of lambda values. Left unspecified, the a sequence of lambda values is automatically computed, ranging uniformly on the log scale over the relevant range of lambda values.
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 parameter estimates 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.
lambda.fused	tuning parameter for fused (generalized) lasso penalty
alpha	currently not used. Will be used later for fused lasso
group.weights	A vector of values representing multiplicative factors by which each group's penalty is to be multiplied. Often, this is a function (such as the square root) of the number of predictors in each group. The default is to use the square root of group size for the group selection methods.
adaptive.lasso	Flag indicating whether or not to use adaptive lasso weights. If set to TRUE and <code>group.weights</code> is unspecified, then this will override <code>group.weights</code> . If a vector is supplied to <code>group.weights</code> , then the <code>adaptive.lasso</code> weights will be multiplied by the <code>group.weights</code> vector
adaptive.fused	Flag indicating whether or not to use adaptive fused lasso weights.
penalty.factor	vector of weights to be multiplied to the tuning parameter for the group lasso penalty. A vector of length equal to the number of groups
penalty.factor.fused	vector of weights to be multiplied to the tuning parameter for the fused lasso penalty. A vector of length equal to the number of variables. mostly for internal usage
gamma	power to raise the MLE estimated weights by for the adaptive lasso. defaults to 1

<code>standardize</code>	Logical flag for x variable standardization, prior to fitting the models. The coefficients are always returned on the original scale. Default is <code>standardize = TRUE</code> . If variables are in the same units already, you might not wish to standardize.
<code>intercept</code>	Should intercept(s) be fitted ( <code>default = TRUE</code> ) or set to zero ( <code>FALSE</code> )
<code>compute.se</code>	Should standard errors be computed? If <code>TRUE</code> , then models are re-fit with no penalization and the standard errors are computed from the refit models. These standard errors are only theoretically valid for the adaptive lasso (when <code>adaptive.lasso</code> is set to <code>TRUE</code> )
<code>rho</code>	ADMM parameter. must be a strictly positive value. By default, an appropriate value is automatically chosen
<code>dynamic.rho</code>	<code>TRUE/FALSE</code> indicating whether or not the rho value should be updated throughout the course of the ADMM iterations
<code>maxit</code>	integer. Maximum number of ADMM iterations. Default is 500.
<code>abs.tol</code>	absolute convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is $10^{-5}$ , which is typically adequate.
<code>rel.tol</code>	relative convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is $10^{-5}$ , which is typically adequate.
<code>irls.tol</code>	convergence tolerance for IRLS iterations. Only used if <code>family != "gaussian"</code> . Default is $10^{-5}$ .
<code>irls.maxit</code>	integer. Maximum number of IRLS iterations. Only used if <code>family != "gaussian"</code> . Default is 100.

**Value**

An object with S3 class "oglasso"

**Examples**

```
library(vennLasso)

set.seed(123)
n.obs <- 1e3
n.vars <- 50

true.beta <- c(rep(0,2), 1, -1, rep(0, 8), 0.5, -0.5, 1, rep(0, 35))

x <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
y <- rnorm(n.obs, sd = 3) + drop(x %*% true.beta)

groups <- c(list(c(1,2), c(2,3), c(3,4,5), 5:10, 6:12, 7:15), lapply(16:50, function(x) x))

## Not run:
fit <- oglasso(x = x, y = y, group = groups)

## End(Not run)
```

---

plot.cv.vennLasso      *Plot method for cv.vennLasso fitted objects*

---

### Description

Plot method for cv.vennLasso fitted objects  
 Plotting method for vennLasso fitted objects

### Usage

```
## S3 method for class 'cv.vennLasso'
plot(x, sign.lambda = 1, ...)

## S3 method for class 'vennLasso'
plot(
  x,
  which.subpop = 1,
  xvar = c("norm", "lambda", "loglambda", "dev"),
  xlab = iname,
  ylab = "Coefficients",
  ...
)
```

### Arguments

x	fitted vennLasso or cv.vennLasso model object
sign.lambda	Either plot against log(lambda) (default) or its negative if sign.lambda = -1.
...	other graphical parameters for the plot
which.subpop	which row in the coefficient matrix should be plotting? Each row corresponds to a particular combination of the specified stratifying variables
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.
xlab	character value supplied for x-axis label
ylab	character value supplied for y-axis label

### Examples

```
set.seed(123)

dat.sim <- genHierSparseData(ncats = 3, nvars = 25,
                             nobs = 100,
                             hier.sparsity.param = 0.5,
                             prop.zero.vars = 0.5,
                             effect.size.max = 0.25,
                             family = "gaussian")
```

```

x      <- dat.sim$x
x.test <- dat.sim$x.test
y      <- dat.sim$y
y.test <- dat.sim$y.test
grp    <- dat.sim$group.ind
grp.test <- dat.sim$group.ind.test

fit.adapt <- cv.vennLasso(x, y,
                        grp,
                        adaptive.lasso = TRUE,
                        nlambda       = 25,
                        nfolds        = 4)

plot(fit.adapt)

library(Matrix)

set.seed(123)
n.obs <- 200
n.vars <- 50

true.beta.mat <- array(NA, dim = c(3, n.vars))
true.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
true.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
true.beta.mat[3,] <- c(0, 0, 1, 1, -1, rep(0, n.vars - 5))
rownames(true.beta.mat) <- c("1,0", "1,1", "0,1")
true.beta <- as.vector(t(true.beta.mat))

x.sub1 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub2 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub3 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)

x <- as.matrix(rbind(x.sub1, x.sub2, x.sub3))

conditions <- as.matrix(cbind(c(rep(1, 2 * n.obs), rep(0, n.obs)),
                             c(rep(0, n.obs), rep(1, 2 * n.obs))))

y <- rnorm(n.obs * 3, sd = 3) + drop(as.matrix(bdiag(x.sub1, x.sub2, x.sub3)) %*% true.beta)

fit <- vennLasso(x = x, y = y, groups = conditions)

layout(matrix(1:3, ncol = 3))
plot(fit, which.subpop = 1)
plot(fit, which.subpop = 2)
plot(fit, which.subpop = 3)

```

**Description**

plotting function to investigate estimated coefficients

**Usage**

```
plotCoefs(object, s = NULL, ...)
```

**Arguments**

object	fitted vennLasso object
s	lambda value for the predictions. Only one can be specified at a time
...	other graphical parameters for the plot

**Examples**

```
set.seed(123)

dat.sim <- genHierSparseData(ncats = 3, nvars = 25, nobs = 200)

fit <- vennLasso(x = dat.sim$x, y = dat.sim$y, groups = dat.sim$group.ind)

plotCoefs(fit, s = fit$lambda[22])
```

---

plotSelections	<i>plotting function to investigate hierarchical structure of selection</i>
----------------	---

---

**Description**

plotting function to investigate hierarchical structure of selection

**Usage**

```
plotSelections(object, s = NULL, type = c("d3.tree"), ...)
```

**Arguments**

object	fitted vennLasso object
s	lambda value for the predictions. Only one can be specified at a time
type	type of plot to make. Currently only "d3.tree" and "igraph.tree" available
...	other graphical parameters for the plot



**Examples**

```
set.seed(123)

dat.sim <- genHierSparseData(ncats = 3, nvars = 25, nobs = 200)

fit <- vennLasso(x = dat.sim$x, y = dat.sim$y, groups = dat.sim$group.ind)

plotSelections(fit, s = fit$lambda[32])
```

---

plotVenn

*plotting function for venn diagrams of overlapping conditions*

---

**Description**

plotting function for venn diagrams of overlapping conditions

**Usage**

```
plotVenn(
  conditions,
  condition.names = NULL,
  lty = "blank",
  fill.colors = c("royalblue1", "goldenrod1", "mediumvioletred", "turquoise3",
    "firebrick1"),
  ...
)
```

**Arguments**

conditions	condition matrix such as the one given to vennLasso() function. It can have up to 5 conditions
condition.names	names of the conditions (equal to the number of columns of conditions)
lty	standard 'lty' graphical parameter for line type around circles. default is no lines
fill.colors	vector of colors for plotting. Set fill.colors = NULL for no colors
...	other graphical parameters for the plot

**Examples**

```
library(Matrix)

set.seed(123)
n.obs <- 200
n.vars <- 50
```

```

true.beta.mat <- array(NA, dim = c(3, n.vars))
true.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
true.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
true.beta.mat[3,] <- c(0, 0, 1, 1, -1, rep(0, n.vars - 5))
rownames(true.beta.mat) <- c("1,0", "1,1", "0,1")
true.beta <- as.vector(t(true.beta.mat))

x.sub1 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub2 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)
x.sub3 <- matrix(rnorm(n.obs * n.vars), n.obs, n.vars)

x <- as.matrix(rbind(x.sub1, x.sub2, x.sub3))

conditions <- as.matrix(cbind(c(rep(1, 2 * n.obs), rep(0, n.obs)),
                             c(rep(0, n.obs), rep(1, 2 * n.obs))))

y <- rnorm(n.obs * 3, sd = 3) + drop(as.matrix(bdiag(x.sub1, x.sub2, x.sub3)) %*% true.beta)

fit <- vennLasso(x = x, y = y, groups = conditions)

vennobj <- plotVenn(conditions)

```

---

predict.cv.vennLasso *Prediction for Cross Validation Hierarchical Lasso Object*

---

## Description

Prediction for Cross Validation Hierarchical Lasso Object

## Usage

```

## S3 method for class 'cv.vennLasso'
predict(object, newx, group.mat, s = c("lambda.min"), use.refit = FALSE, ...)

```

## Arguments

object	fitted cv.vennLasso object
newx	new matrix for predictions
group.mat	A matrix of the group memberships for now. Ignore the rest: A list of length equal to the number of groups containing vectors of integers indicating the variable IDs for each group. For example, groups=list(c(1,2), c(2,3), c(3,4,5)) specifies that Group 1 contains variables 1 and 2, Group 2 contains variables 2 and 3, and Group 3 contains variables 3, 4, and 5. Can also be a matrix of 0s and 1s with the number of columns equal to the number of groups and the number of rows equal to the number of variables. A value of 1 in row i and column j indicates that variable i is in group j and 0 indicates that variable i is not in group j.

s	lambda value for the predictions. defaults to all values computed in the vennLasso object
use.refit	Should the refitted beta estimates be used for prediction? Defaults to FALSE. If TRUE then the beta estimates from the model refit on just the selected covariates are used
...	parameters to be passed to predict.vennLasso

**Value**

predictions or coefficients

---

predict.vennLasso      *Prediction for Hierarchical Shared Lasso*

---

**Description**

Prediction for Hierarchical Shared Lasso

**Usage**

```
## S3 method for class 'vennLasso'
predict(
  object,
  newx,
  group.mat,
  s = NULL,
  use.refit = FALSE,
  type = c("link", "response", "coefficients", "nonzero", "class", "nvars", "median",
           "survival"),
  ...
)
```

**Arguments**

object	fitted vennLasso object
newx	new matrix for predictions
group.mat	A matrix of the group memberships for now. Ignore the rest: A list of length equal to the number of groups containing vectors of integers indicating the variable IDs for each group. For example, groups=list(c(1,2), c(2,3), c(3,4,5)) specifies that Group 1 contains variables 1 and 2, Group 2 contains variables 2 and 3, and Group 3 contains variables 3, 4, and 5. Can also be a matrix of 0s and 1s with the number of columns equal to the number of groups and the number of rows equal to the number of variables. A value of 1 in row i and column j indicates that variable i is in group j and 0 indicates that variable i is not in group j.

s	lambda value for the predictions. defaults to all values computed in the vennLasso object
use.refit	Should the refitted beta estimates be used for prediction? Defaults to FALSE. If TRUE then the beta estimates from the model refit on just the selected covariates are used
type	type of predictions to be made. type = "median" is for the median survival time and type = "survival" is for the predicted hazard function
...	parameters to be passed to vennLasso

**Value**

predictions or coefficients

---

vennLasso	<i>Fitting vennLasso models</i>
-----------	---------------------------------

---

**Description**

Fitting vennLasso models

**Usage**

```
vennLasso(
  x,
  y,
  groups,
  family = c("gaussian", "binomial"),
  nlambda = 100L,
  lambda = NULL,
  lambda.min.ratio = NULL,
  lambda.fused = NULL,
  penalty.factor = NULL,
  group.weights = NULL,
  adaptive.lasso = FALSE,
  adaptive.fused = FALSE,
  gamma = 1,
  standardize = FALSE,
  intercept = TRUE,
  one.intercept = FALSE,
  compute.se = FALSE,
  conf.int = NULL,
  rho = NULL,
  dynamic.rho = TRUE,
  maxit = 500L,
  abs.tol = 1e-05,
  rel.tol = 1e-05,
```

```

  irls.tol = 1e-05,
  irls.maxit = 100L,
  model.matrix = FALSE,
  ...
)

```

## Arguments

<code>x</code>	input matrix of dimension <code>nobs</code> by <code>nvars</code> . Each row is an observation, each column corresponds to a covariate
<code>y</code>	numeric response vector of length <code>nobs</code>
<code>groups</code>	A list of length equal to the number of groups containing vectors of integers indicating the variable IDs for each group. For example, <code>groups = list(c(1, 2), c(2, 3), c(3, 4, 5))</code> specifies that Group 1 contains variables 1 and 2, Group 2 contains variables 2 and 3, and Group 3 contains variables 3, 4, and 5. Can also be a matrix of 0s and 1s with the number of columns equal to the number of groups and the number of rows equal to the number of variables. A value of 1 in row <i>i</i> and column <i>j</i> indicates that variable <i>i</i> is in group <i>j</i> and 0 indicates that variable <i>i</i> is not in group <i>j</i> .
<code>family</code>	"gaussian" for least squares problems, "binomial" for binary response, and "coxph" for time-to-event outcomes (not yet available)
<code>nlambda</code>	The number of lambda values. Default is 100.
<code>lambda</code>	A user-specified sequence of lambda values. Left unspecified, the a sequence of lambda values is automatically computed, ranging uniformly on the log scale over the relevant range of lambda values.
<code>lambda.min.ratio</code>	Smallest value for lambda, as a fraction of <code>lambda.max</code> , the (data derived) entry value (i.e. the smallest value for which all parameter estimates 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> can lead to a saturated fit when <code>nobs &lt; nvars</code> .
<code>lambda.fused</code>	tuning parameter for fused lasso penalty
<code>penalty.factor</code>	vector of weights to be multiplied to the tuning parameter for the group lasso penalty. A vector of length equal to the number of groups
<code>group.weights</code>	A vector of values representing multiplicative factors by which each group's penalty is to be multiplied. Often, this is a function (such as the square root) of the number of predictors in each group. The default is to use the square root of group size for the group selection methods.
<code>adaptive.lasso</code>	Flag indicating whether or not to use adaptive lasso weights. If set to TRUE and <code>group.weights</code> is unspecified, then this will override <code>group.weights</code> . If a vector is supplied to <code>group.weights</code> , then the adaptive.lasso weights will be multiplied by the <code>group.weights</code> vector
<code>adaptive.fused</code>	Flag indicating whether or not to use adaptive fused lasso weights.
<code>gamma</code>	power to raise the MLE estimated weights by for the adaptive lasso. defaults to 1

<code>standardize</code>	Should the data be standardized? Defaults to FALSE.
<code>intercept</code>	Should an intercept be fit? Defaults to TRUE
<code>one.intercept</code>	Should a single intercept be fit for all subpopulations instead of one for each subpopulation? Defaults to FALSE.
<code>compute.se</code>	Should standard errors be computed? If TRUE, then models are re-fit with no penalization and the standard errors are computed from the refit models. These standard errors are only theoretically valid for the adaptive lasso (when <code>adaptive.lasso</code> is set to TRUE)
<code>conf.int</code>	level for confidence intervals. Defaults to NULL (no confidence intervals). Should be a value between 0 and 1. If confidence intervals are to be computed, <code>compute.se</code> will be automatically set to TRUE
<code>rho</code>	ADMM parameter. must be a strictly positive value. By default, an appropriate value is automatically chosen
<code>dynamic.rho</code>	TRUE/FALSE indicating whether or not the rho value should be updated throughout the course of the ADMM iterations
<code>maxit</code>	integer. Maximum number of ADMM iterations. Default is 500.
<code>abs.tol</code>	absolute convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is $10^{-5}$ , which is typically adequate.
<code>rel.tol</code>	relative convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is $10^{-5}$ , which is typically adequate.
<code>irls.tol</code>	convergence tolerance for IRLS iterations. Only used if <code>family != "gaussian"</code> . Default is $10^{-5}$ .
<code>irls.maxit</code>	integer. Maximum number of IRLS iterations. Only used if <code>family != "gaussian"</code> . Default is 100.
<code>model.matrix</code>	logical flag. Should the design matrix used be returned?
<code>...</code>	not used

**Value**

An object with S3 class "vennLasso"

**Examples**

```
library(Matrix)

# first simulate heterogeneous data using
# genHierSparseData
set.seed(123)
dat.sim <- genHierSparseData(ncats = 2, nvars = 25,
                             nobs = 200,
                             hier.sparsity.param = 0.5,
                             prop.zero.vars = 0.5,
                             family = "gaussian")

x <- dat.sim$x
conditions <- dat.sim$group.ind
```

```

y          <- dat.sim$y

true.beta.mat <- dat.sim$beta.mat

fit <- vennLasso(x = x, y = y, groups = conditions)

(true.coef <- true.beta.mat[match(dimnames(fit$beta)[[1]], rownames(true.beta.mat)),])
round(fit$beta[, ,21], 2)

## fit adaptive version and compute confidence intervals
afit <- vennLasso(x = x, y = y, groups = conditions, conf.int = 0.95, adaptive.lasso = TRUE)

(true.coef <- true.beta.mat[match(dimnames(fit$beta)[[1]], rownames(true.beta.mat)),,][,1:10]
round(afit$beta[,1:10,28], 2)
round(afit$lower.ci[,1:10,28], 2)
round(afit$upper.ci[,1:10,28], 2)

aic.idx <- which.min(afit$aic)
bic.idx <- which.min(afit$bic)

# actual coverage
# actual coverage
mean(true.coef[afit$beta[, -1, aic.idx] != 0] >=
      afit$lower.ci[, -1, aic.idx][afit$beta[, -1, aic.idx] != 0] &
      true.coef[afit$beta[, -1, aic.idx] != 0] <=
      afit$upper.ci[, -1, aic.idx][afit$beta[, -1, aic.idx] != 0])

(covered <- true.coef >= afit$lower.ci[, -1, aic.idx] & true.coef <= afit$upper.ci[, -1, aic.idx])
mean(covered)

# logistic regression example
## Not run:
set.seed(123)
dat.sim <- genHierSparseData(ncats = 2, nvars = 25,
                            nobs = 200,
                            hier.sparsity.param = 0.5,
                            prop.zero.vars = 0.5,
                            family = "binomial",
                            effect.size.max = 0.5) # don't make any
                                                    # coefficients too big

x          <- dat.sim$x
conditions <- dat.sim$group.ind
y          <- dat.sim$y
true.beta.b <- dat.sim$beta.mat

bfit <- vennLasso(x = x, y = y, groups = conditions, family = "binomial")

(true.coef.b <- -true.beta.b[match(dimnames(fit$beta)[[1]], rownames(true.beta.b)),])
round(bfit$beta[, ,20], 2)

## End(Not run)

```





# Index

`cv.vennLasso`, [2](#)

`estimate.hier.sparsity.param`, [5](#)

`genHierSparseBeta`, [6](#)

`genHierSparseData`, [7](#)

`logLik.vennLasso`, [10](#)

`oglasso`, [11](#)

`plot.cv.vennLasso`, [14](#)

`plot.vennLasso (plot.cv.vennLasso)`, [14](#)

`plotCoefs`, [15](#)

`plotSelections`, [16](#)

`plotVenn`, [17](#)

`predict.cv.vennLasso`, [18](#)

`predict.vennLasso`, [19](#)

`vennLasso`, [20](#)