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>.

**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

**RooxygenNote**  7.1.0

**Suggests**  knitr, rmarkdown

**VignetteBuilder**  knitr

**NeedsCompilation**  yes

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**Repository**  CRAN

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Cross Validation for the vennLasso

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 glmnet, 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 foldid 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 doMC.

... parameters to be passed to vennLasso

Value
An object with S3 class "cv.vennLasso"
library(Matrix)
set.seed(123)
n.obs <- 150
n.vars <- 25

ture.beta.mat <- array(NA, dim = c(3, n.vars))
ture.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
ture.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
ture.beta.mat[3,] <- c(0, 0, 1, 1, -1, rep(0, n.vars - 5))
rownames(ture.beta.mat) <- c("1,0", "1,1", "0,1")
ture.beta <- as.vector(t(ture.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)) %*% ture.beta)
fit <- cv.vennLasso(x = x, y = y, groups = conditions, nfolds = 3)
fitted.coef <- predict(fit$vennLasso.fit, type = "coefficients", s = fit$lambda.min)
(ture.coef <- ture.beta.mat[match(dimnames(fit$vennLasso.fit$beta)[[1]],
                                 rownames(ture.beta.mat))],)
round(fitted.coef, 2)
## effects smaller for logistic regression
Not run:
ture.beta.mat <- ture.beta.mat / 2
ture.beta <- ture.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)) %*% ture.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)
(ture.coef <- ture.beta.mat[match(dimnames(bfit$vennLasso.fit$beta)[[1]],
                                 rownames(ture.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

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**Description**

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

**Usage**

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

**Arguments**

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

**Examples**

```r
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
```
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,  nob,
    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

ncats number of categories to stratify on
nvars number of variables
nobs number of observations per strata to simulate
nobs.test number of independent test observations per strata to simulate
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

prop.zero.vars proportion of all variables that will be zero across all strata
effect.size.max maximum magnitude of the true effect sizes
misspecification.prop proportion of variables with hierarchical missingness misspecified
family family for the response variable
sd standard deviation for gaussian simulations
snr signal-to-noise ratio (only used for family = "gaussian")
beta a matrix of true beta values. If given, then no beta will be created and data will be simulated from the given beta
tau rate parameter for rexp() for generating time-to-event outcomes
covar 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:
hs <- 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,
nlambda = 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

ttrue.beta.mat <- array(NA, dim = c(3, n.vars))
ttrue.beta.mat[1,] <- c(-0.5, -1, 0, 0, 2, rep(0, n.vars - 5))
ttrue.beta.mat[2,] <- c(0.5, 0.5, -0.5, -0.5, 1, -1, rep(0, n.vars - 6))
ttrue.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")
ttrue.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)),
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))) %*% ttrue.beta

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

logLik(fit)
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, `group = 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.

**fused**

Matrix specifying generalized lasso penalty formulation. Each column corresponds to each variable and each row corresponds to a new penalty term, i.e., if row 1 has the first entry of 1 and the second entry of -1, then the penalty term `lambda.fused * |beta_1 - beta_2|` 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, 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 `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.

**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 `group.weights` is unspecified, then this will override `group.weights`. If a vector is supplied to `group.weights`, then the adaptive lasso weights will be multiplied by the `group.weights` 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.
standardize Logical flag for x variable standardization, prior to fitting the models. The coefficients are always returned on the original scale. Default is standardize = TRUE. If variables are in the same units already, you might not wish to standardize.

intercept Should intercept(s) be fitted (default = TRUE) or set to zero (FALSE)

compute.se 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 adaptive.lasso is set to TRUE)

rho ADMM parameter. must be a strictly positive value. By default, an appropriate value is automatically chosen

dynamic.rho TRUE/FALSE indicating whether or not the rho value should be updated throughout the course of the ADMM iterations

maxit integer. Maximum number of ADMM iterations. Default is 500.

abs.tol absolute convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is 10^-5, which is typically adequate.

rel.tol relative convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is 10^-5, which is typically adequate.

irls.tol convergence tolerance for IRLS iterations. Only used if family != "gaussian". Default is 10^-5.

irls.maxit integer. Maximum number of IRLS iterations. Only used if family != "gaussian". 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)
Description
Plot method for cv.vennLasso fitted objects
Plotting method for vennLasso fitted objects

Usage

```r
## 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

```r
set.seed(123)

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

---

plotCoefs

plotting function to investigate estimated coefficients
plotSelections

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
... other graphical parameters for the plot
Examples

set.seed(123)

dat.sim <- genHierSparseData(ncats = 3, nvars = 25, nob = 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, 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**

```r
## 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.
predict.vennLasso

### Description

Prediction for Hierarchical Shared Lasso

### Usage

```r
## 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

### Value

predictions or coefficients
vennLasso

Fitting vennLasso models

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

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

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.

family "gaussian" for least squares problems, "binomial" for binary response, and "coxph" for time-to-event outcomes (not yet available)

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 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 can lead to a saturated fit when nobs < nvars.

lambda.fused tuning parameter for fused lasso penalty

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

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 group.weights is unspecified, then this will override group.weights. If a vector is supplied to group.weights, then the adaptive.lasso weights will be multiplied by the group.weights vector

adaptive.fused Flag indicating whether or not to use adaptive fused lasso weights.

gamma power to raise the MLE estimated weights by for the adaptive lasso. defaults to 1
standardize Should the data be standardized? Defaults to FALSE.
intercept Should an intercept be fit? Defaults to TRUE
one.intercept Should a single intercept be fit for all subpopulations instead of one for each subpopulation? Defaults to FALSE.
compute.se 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 adaptive.lasso is set to TRUE)
conf.int 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, compute.se will be automatically set to TRUE
rho ADMM parameter. must be a strictly positive value. By default, an appropriate value is automatically chosen
dynamic.rho TRUE/FALSE indicating whether or not the rho value should be updated throughout the course of the ADMM iterations
maxit integer. Maximum number of ADMM iterations. Default is 500.
abs.tol absolute convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is 10^{-5}, which is typically adequate.
rel.tol relative convergence tolerance for ADMM iterations for the relative dual and primal residuals. Default is 10^{-5}, which is typically adequate.
irls.tol convergence tolerance for IRLS iterations. Only used if family != "gaussian". Default is 10^{-5}.
irls.maxit integer. Maximum number of IRLS iterations. Only used if family != "gaussian". Default is 100.
model.matrix logical flag. Should the design matrix used be returned?
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

... not used
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(afit$beta)[[1]], rownames(true.beta.mat)),][,1:10])
round(afit$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
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$x
categories <- dat.sim$group.ind
y <- dat.sim$y
true.beta.b <- dat.sim$beta.mat

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

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

## End(Not run)
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