Package ‘accSDA’

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Title Accelerated Sparse Discriminant Analysis

Imports MASS (>= 7.3.45), rARPACK (>= 0.10.0), sparseLDA (>= 0.1.7),
   ggplot2 (>= 2.1.0), ggrepthemes (>= 3.2.0), grid (>= 3.2.2),
   gridExtra (>= 2.2.1)

Depends R (>= 3.2)

Description Implementation of sparse linear discriminant analysis, which is a supervised
classification method for multiple classes. Various novel optimization approaches to
this problem are implemented including alternating direction method of multipliers (ADMM),
proximal gradient (PG) and accelerated proximal gradi-
ent (APG) (See Atkins et al. <arXiv:1705.07194>).
Functions for performing cross validation are also supplied along with basic prediction
and plotting functions.
Sparse zero variance discriminant analysis (SZVD) is also included in the package
(See Ames and Hong. <arXiv:1401.5492>). See the github wiki for a more extended description.

License GPL (>= 2)

URL https://github.com/gumeo/accSDA/wiki

BugReports https://github.com/gumeo/accSDA/issues

LazyData TRUE

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R topics documented:

- accSDA: A package for performing sparse discriminant analysis in various ways.
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Description

The accSDA package provides functions to perform sparse discriminant analysis using a selection of three optimization methods, proximal gradient (PG), accelerated proximal gradient (APG) and alternating direction method of multipliers (ADMM). The package is intended to extend the available tools to perform sparse discriminant analysis in R. The three methods can be called from the function `asda`. Cross validation is also implemented for the L1 regularization parameter. Functions for doing predictions, summary, printing and simple plotting are also provided. The sparse discriminant functions perform lda on the projected data by default, using the lda function in the MASS package. The functions return an object of the same class as the name of the function and provide the lda solution, along with the projected data, thus other kinds of classification algorithms can be employed on the projected data.

Description

Applies accelerated proximal gradient algorithm, proximal gradient algorithm or alternating direction methods of multipliers algorithm to the optimal scoring formulation of sparse discriminant analysis proposed by Clemmensen et al. 2011.

\[
\arg\min_{\beta} |Y_t \theta - X_t \beta|^2_2 + t|\beta|_1 + \lambda|\beta|^2_2
\]
Usage

ASDA(Xt, ...)

## Default S3 method:
ASDA(Xt, Yt, Om = diag(p), gam = 0.001, lam = 1e-06, 
q = K - 1, method = "SDAAP", control = list(), ...)

Arguments

Xt n by p data matrix, (can also be a data.frame that can be coerced to a matrix)

... Additional arguments for `lda` function in package MASS.

Yt n by K matrix of indicator variables (Yij = 1 if i in class j). This will later be 
changed to handle factor variables as well. Each observation belongs in a single 
class, so for a given row/observation, only one element is 1 and the rest is 0.

Om p by p parameter matrix Omega in generalized elastic net penalty.

gam Regularization parameter for elastic net penalty.

lam Regularization parameter for l1 penalty, must be greater than zero. If cross-
validation is used (`CV = TRUE`) then this must be a vector of length greater than 
one.

q Desired number of discriminant vectors.

method This parameter selects which optimization method to use. It is specified as a 
character vector which can be one of the three values

SDAP Proximal gradient algorithm.

SDAAP Accelerated proximal gradient algorithm.

SDAD Alternating directions method of multipliers algorithm.

ccontrol List of control arguments. See Details.

Details

The control list contains the following entries to further tune the algorithms.

PGsteps Maximum number if inner proximal gradient/ADMM algorithm for finding beta. Default 
value is 1000.

PGtol Stopping tolerance for inner method. If the method is SDAD, then this must be a vector of 
two values, absolute (first element) and relative tolerance (second element). Default value is 
1e-5 for both absolute and relative tolerances.

maxits Number of iterations to run. Default value is 250.

tol Stopping tolerance. Default value is 1e-3.

mu Penalty parameter for augmented Lagrangian term, must be greater than zero and only needs to 
be specified when using method SDAD. Default value is 1.

CV Logical value which is TRUE if cross validation is supposed to be performed. If cross-validation 
is performed, then lam should be specified as a vector containing the regularization values to 
be tested. Default value is FALSE.
folds  Integer determining the number of folds in cross-validation. Not needed if CV is not specified. Default value is 5.
feat  Maximum fraction of nonzero features desired in validation scheme. Not needed if CV is not specified. Default value is 0.15.
quiet  Set to FALSE if status updates are supposed to be printed to the R console. Default value is TRUE. Note that this triggers a lot of printing to the console.

Value

ASDA returns an object of class "ASDA" including a list with the following named components:

call  The matched call.
B  p by q matrix of discriminant vectors, i.e. sparse loadings.
Q  K by q matrix of scoring vectors, i.e. optimal scores.
varNames  Names of the predictors used, i.e. column names of Xt.
origP  Number of variables in Xt.
fit  Output from function lda on projected data. This is NULL the trivial solution is found, i.e. B is all zeroes. Use lower values of lam if that is the case.
classes  The classes in Yt.
lambda  The lambda/lam used, best value found by cross-validation if CV is TRUE.

NULL

Note

The input matrix Xt should be normalized, i.e. each column corresponding to a variable should have its mean subtracted and scaled to unit length. The functions normalize and normalizetest are supplied for this purpose in the package.

See Also

SDAAP, SDAP and SDAD

Examples

set.seed(123)
# Prepare training and test set
train <- c(1:40,51:90,101:140)
Xtrain <- iris[train,1:4]
nX <- normalize(Xtrain)
Xtrain <- nX$x
Ytrain <- iris[train,5]
Xtest <- iris[-train,1:4]
Xtest <- normalizetest(Xtest,nX)
Ytest <- iris[-train,5]

# Define parameters for Alternating Direction Method of Multipliers (SDAD)
Om <- diag(4)+0.1*matrix(1,4,4) #elNet coef mat
gam <- 0.01
lam <- 0.01
method <- "SDAD"
q <- 2
control <- list(PGsteps = 100,
                PGtol = c(1e-5, 1e-5),
                mu = 1,
                maxits = 100,
                tol = 1e-3,
                quiet = FALSE)

# Run the algorithm
res <- ASDA(Xt = Xtrain,
            Yt = Ytrain,
            Om = Om,
            gam = gam,
            lam = lam,
            q = q,
            method = method,
            control = control)

# Can also just use the defaults, which is Accelerated Proximal Gradient (SDAAP):
resDef <- ASDA(Xtrain,Ytrain)

# Some example on simulated data
# Generate Gaussian data on three classes with plenty of redundant variables

# This example shows the basic steps on how to apply this to data, i.e.:
# 1) Setup training data
# 2) Normalize
# 3) Train
# 4) Predict
# 5) Plot projected data
# 6) Accuracy on test set

P <- 300 # Number of variables
N <- 50 # Number of samples per class

# Mean for classes, they are zero everywhere except the first 3 coordinates
m1 <- rep(0,P)
m1[1] <- 3

m2 <- rep(0,P)
m2[2] <- 3

m3 <- rep(0,P)
m3[3] <- 3

# Sample dummy data
Xtrain <- rbind(MASS::mvrnorm(n=N,mu = m1, Sigma = diag(P)),
                MASS::mvrnorm(n=N,mu = m2, Sigma = diag(P)),
                MASS::mvrnorm(n=N,mu = m3, Sigma = diag(P)))
Xtest <- rbind(MASS::mvrnorm(n=N,mu = m1, Sigma = diag(P)),
       MASS::mvrnorm(n=N,mu = m2, Sigma = diag(P)),
       MASS::mvrnorm(n=N,mu = m3, Sigma = diag(P)))

# Generate the labels
Ytrain <- factor(rep(1:3,each=N))
Ytest <- Ytrain

# Normalize the data
Xt <- accSDA::normalize(Xtrain)
Xtrain <- Xt$Xc # Use the centered and scaled data
Xtest <- accSDA::normalizetest(Xtest,Xt)

# Train the classifier and increase the sparsity parameter from the default
# so we penalize more for non-sparse solutions.
res <- accSDA::ASDA(Xtrain,Ytrain,lambda=0.01)

# Plot the projected training data, it is projected to
# 2-dimension because we have 3 classes. The number of discriminant
# vectors is maximum number of classes minus 1.
XtrainProjected <- Xtrain%*%res$beta
plot(XtrainProjected[,1],XtrainProjected[,2],col=Ytrain)

# Predict on the test data
preds <- predict(res, newdata = Xtest)

# Plot projected test data with predicted and correct labels
XtestProjected <- Xtest%*%res$beta
plot(XtestProjected[,1],XtestProjected[,2],col=Ytest,
     main="Projected test data with original labels")
plot(XtestProjected[,1],XtestProjected[,2],col=preds$class,
     main="Projected test data with predicted labels")

# Calculate accuracy
sum(preds$class == Ytest)/(3*N) # We have N samples per class, so total 3*N

---

**ASDABarPlot**

*barplot for ASDA objects*

**Description**

This is a function to visualize the discriminant vector from the ASDA method. The plot is constructed as a ggplot barplot and the main purpose of it is to visually inspect the sparsity of the discriminant vectors. The main things to look for are how many parameters are non-zero and if there is any structure in the ones that are non-zero, but the structure is dependent on the order you specify your variables. For time-series data, this could mean that a chunk of variables are non-zero that are close in time, meaning that there is some particular event that is best for discriminating between the classes that you have.
ASDABarPlot

Usage

ASDABarPlot(asdaObj, numDVs = 1, xlabel, ylabel, getList = FALSE, main, ...)

Arguments

asdaObj Object from the ASDA function.
numDVs Number of discriminant vectors (DVs) to plot. This is limited by the number of DVs outputted from the ASDA function or k-1 DVs where k is the number of classes. The first 1 to numDVs are plotted.
xlabel Label to put under every plot
ylabel Vector of y-axis labels for each plot, e.g. if there are three DVs, then ylab = c('Discriminant Vector 1', 'Discriminant Vector 2', 'Discriminant Vector 3') is a valid option.
geList Logical value indicating whether the output should be a list of the plots or the plots stacked in one plot using the gridExtra package. By default the function produces a single plot combining all plots of the DVs.
main Main title for the plots, this is not used if getList is set to TRUE.
... Extra arguments to grid.arrange.

Value

barplot.ASDA returns either a single combined plot or a list of individual ggplot objects.

Note

This function is used as a quick diagnostics tool for the output from the ASDA function. Feel free to look at the code to customize the plots in any way you like.

See Also

ASDA

Examples

# Generate and ASDA object with your data, e.g.
# Prepare training and test set
# This is a very small data set, I advise you to try it on something with more variables, e.g. something from this source: http://www.cs.ucr.edu/~eamonn/time_series_data/
# or possibly run this on the Gaussian data example from the ASDA function
train <- c(1:40,51:90,101:140)
Xtrain <- iris[train,1:4]
nX <- normalize(Xtrain)
Xtrain <- nX$Xc
Ytrain <- iris[train,5]
Xtest <- iris[-train,1:4]
Xtest <- normalize(Xtest,nX)
Ytest <- iris[-train,5]
# Run the method
resIris <- ASDA(Xtrain,Ytrain)
# Look at the barplots of the DVs
ASDABarPlot(resIris)

normalize

**Normalize training data**

**Description**
Normalize a vector or matrix to zero mean and unit length columns.

**Usage**
normalize(X)

**Arguments**
- **X**: a matrix with the training data with observations down the rows and variables in the columns.

**Details**
This function can e.g. be used for the training data in the ASDA function.

**Value**
normalize Returns a list with the following attributes:
- **Xc**: The normalized data
- **mx**: Mean of columns of X.
- **vx**: Length of columns of X.
- **Id**: Logical vector indicating which variables are included in X. If some of the columns have zero length they are omitted.

**Author(s)**
Line Clemmensen

**References**

**See Also**
ormalizetest, predict.ASDA, ASDA
normalizetest

Examples

```r
## Data
X <- matrix(sample(seq(3), 12, replace = TRUE), nrow = 3)

## Normalize data
Nm <- normalize(X)
print(Nm$Xc)

## See if any variables have been removed
which(!Nm$Id)
```

normalizetest  Normalize training data

Description

Normalize test data using output from the `normalize()` of the training data

Usage

`normalizetest(Xtst, Xn)`

Arguments

- `Xtst`: a matrix with the test data with observations down the rows and variables in the columns.
- `Xn`: List with the output from `normalize(Xtr)` of the training data.

Details

This function can e.g. be used for the test data in the `predict.ASDA` function.

Value

`normalizetest` returns the normalized test data `Xtst`

Author(s)

Line Clemmensen

References


See Also

`normalize`, `predict.ASDA`, `ASDA`
Examples

```r
## Data
Xtr <- matrix(sample(seq(3), 12, replace=TRUE), nrow=3)
Xtst <- matrix(sample(seq(3), 12, replace=TRUE), nrow=3)

## Normalize training data
Nm <- normalize(Xtr)

## Normalize test data
Xtst <- normalizetest(Xtst, Nm)
```

predict.ASDA

*Predict method for sparse discriminant analysis*

Description

Predicted values based on fit from the function `ASDA`. This function is used to classify new observations based on their explanatory variables/features.

Usage

```r
## S3 method for class 'ASDA'
predict(object, newdata = NULL, ...)
```

Arguments

- **object**
  - Object of class ASDA. This object is returned from the function `ASDA`.
- **newdata**
  - A matrix of new observations to classify.
- **...**
  - Arguments passed to `predict.lda`.

Value

A list with components:

- **class**
  - The classification (a factor)
- **posterior**
  - posterior probabilities for the classes
- **x**
  - the scores

Note

The input matrix `newdata` should be normalized w.r.t. the normalization of the training data.

See Also

`SDAAP`, `SDAP` and `SDAD`
Examples

# Prepare training and test set
train <- c(1:40, 51:90, 101:140)
Xtrain <- iris[train, 1:4]
nX <- normalize(Xtrain)
Xtrain <- nX$Xc
Ytrain <- iris[train, 5]
Xtest <- iris[-train, 1:4]
Xtest <- normalize(Xtest, nX)
Ytest <- iris[-train, 5]

# Define parameters for SDAD
Om <- diag(4)*0.1*matrix(1,4,4) #elNet coef mat
gam <- 0.01
lam <- 0.01
method <- "SDAD"
q <- 2
control <- list(PGsteps = 100,
                PGtol = c(1e-5, 1e-5),
                mu = 1,
                maxits = 100,
                tol = 1e-3,
                quiet = FALSE)

# Run the algorithm
res <- ASDA(Xt = Xtrain,
            Yt = Ytrain,
            Om = Om,
            gam = gam,
            lam = lam,
            q = q,
            method = method,
            control = control)

# Do the predictions on the test set
preds <- predict(object = res, newdata = Xtest)

print.ASDA

Print method for ASDA object

Description

Prints a summary of the output from the ASDA function. The output summarizes the discriminant analysis in human readable format.

Usage

## S3 method for class 'ASDA'
print(x, digits = max(3,getOption("digits") - 3),
      numshow = 5, ...)
Arguments

- **x**: Object of class ASDA. This object is returned from the function `ASDA`.
- **digits**: Number of digits to show in printed numbers.
- **numshow**: Number of best ranked variables w.r.t. to their absolute coefficients.
- **...**: arguments passed to or from other methods.

Value

An invisible copy of `x`.

See Also

- `ASDA`
- `predict.ASDA`
- `SDAD`

Examples

```r
# Prepare training and test set
train <- c(1:40, 51:90, 101:140)
Xtrain <- iris[train, 1:4]
NX <- normalize(Xtrain)
Xtrain <- NX$Xc
Ytrain <- iris[train, 5]
Xtest <- iris[-train, 1:4]
Xtest <- normalize(test(Xtest, NX)
Ytest <- iris[-train, 5]

# Run the algorithm
resDef <- ASDA(Xtrain, Ytrain)

# Print
print(resDef)
```

---

**SZVD**

*Sparse Zero Variance Discriminant Analysis*

Description

Applies SZVD heuristic for sparse zero-variance discriminant analysis to given training set.

Usage

```r
SZVD(train, ...)
```

## Default S3 method:

```r
SZVD(train, gamma, D, penalty = TRUE, scaling = TRUE,
     tol = list(abs = 1e-04, rel = 1e-04), maxits = 2000, beta = 1,
     quiet = TRUE)
```
Arguments

- **train**: Data matrix where first column is the response class.
- **...**: Parameters passed to SZVD.default.
- **gamma**: Set of regularization parameters controlling l1-penalty.
- **D**: dictionary/basis matrix.
- **penalty**: Controls whether to apply reweighting of l1-penalty (using sigma = within-class std devs).
- **scaling**: Logical indicating whether to scale data such that each feature has variance 1.
- **tol**: Stopping tolerances for ADMM algorithm, must include tol$rel and tol$sabs.
- **maxits**: Maximum number of iterations used in the ADMM algorithm.
- **beta**: penalty term controlling the splitting constraint.
- **quiet**: Print intermediate output or not.

Details

This function will currently solve as a standalone function in accSDA for time comparison. A wrapper function like ASDA will be created to use the functionality of plots and such. Maybe call it ASZDA. For that purpose the individual ZVD function will need to be implemented.

Value

**SZVD** returns an object of class "SZVD" including a list with the following named components:

- **DVs**: Discriminant vectors.
- **its**: Number of iterations required to find DVs.
- **pen_scal**: Weights used in reweighted l1-penalty.
- **N**: Basis for the null-space of the sample within-class covariance.
- **means**: Training class-means.
- **mus**: Training mean and variance scaling/centering terms.
- **w0**: unpenalized zero-variance discriminants (initial solutions) plus B and W, etc.

See Also

Used by: **SZVDcv**.

Examples

```r
set.seed(123)
P <- 300 # Number of variables
N <- 50 # Number of samples per class

# Mean for classes, they are zero everywhere except the first 3 coordinates
m1 <- rep(0,P)
```
m1[1] <- 3
m2 <- rep(0,P)
m2[2] <- 3
m3 <- rep(0,P)
m3[3] <- 3

# Sample dummy data
Xtrain <- rbind(MASS::mvrnorm(n=N,mu = m1, Sigma = diag(P)),
               MASS::mvrnorm(n=N,mu = m2, Sigma = diag(P)),
               MASS::mvrnorm(n=N,mu = m3, Sigma = diag(P)))

# Generate the labels
Ytrain <- rep(1:3,each=N)

# Normalize the data
Xt <- accSDA::normalize(Xtrain)
Xtrain <- Xt%*%Xc

# Train the classifier and increase the sparsity parameter from the default
# so we penalize more for non-sparse solutions.
res <- accSDA::SZVD(cbind(Ytrain,Xtrain),beta=2.5,
                     maxits=1000,tol = list(abs = 1e-04, rel = 1e-04))

---

**SZVDcv**

*Cross-validation of sparse zero variance discriminant analysis*

**Description**

Applies alternating direction methods of multipliers to solve sparse zero variance discriminant analysis.

**Usage**

```r
SZVDcv(Atrain, ...)
```

## Default S3 method:

```r
SZVDcv(Atrain, Aval, k, num_gammas, g_mults, D, sparsity_pen,
       scaling, penalty, beta, tol, ztol, maxits, quiet)
```

**Arguments**

- **Atrain**
  - Training data set.
- **...**
  - Parameters passed to `SZVD.default`.
- **Aval**
  - Validation set.
- **k**
  - Number of classes within training and validation sets.
num_gammas  Number of gammas to train on.
g_mults  Parameters defining range of gammas to train, g_max*(c_min, c_max). Note that it is an array/vector with two elements.
D  Penalty dictionary basis matrix.
sparsity_pen  weight defining validation criteria as weighted sum of misclassification error and cardinality of discriminant vectors.
scaling  Whether to rescale data so each feature has variance 1.
penalty  Controls whether to apply reweighting of l1-penalty (using sigma = within-class std devs)
beta  Parameter for augmented Lagrangian term in the ADMM algorithm.
tol  Stopping tolerances for the ADMM algorithm, must have tol$rel and tol$abs.
ztol  Threshold for truncating values in DVs to zero.
maxits  Maximum number of iterations used in the ADMM algorithm.
quiet  Controls display of intermediate results.

Details

This function might require a wrapper similar to ASDA.

Value

SZVDcv returns an object of class "SZVDcv" including a list with the following named components:

DV's  Discriminant vectors for the best choice of gamma.
all_DV's  Discriminant vectors for all choices of gamma.
10_DV's  Discriminant vectors for gamma minimizing cardinality.
mc_DV's  Discriminant vector minimizing misclassification.
gamma  Choice of gamma minimizing validation criterion.
gammas  Set of all gammas trained on.
max_g  Maximum value of gamma guaranteed to yield a nontrivial solution.
ind  Index of best gamma.
w0  unpenalized zero-variance discriminants (initial solutions) plus B and W, etc. from ZVD

See Also

Non CV version: SZVD.
Examples

\begin{verbatim}
P <- 300 # Number of variables
N <- 50 # Number of samples per class

# Mean for classes, they are zero everywhere except the first 3 coordinates
m1 <- rep(0,P)
m1[1] <- 3
m2 <- rep(0,P)
m2[2] <- 3
m3 <- rep(0,P)
m3[3] <- 3

# Sample dummy data
Xtrain <- rbind(MASS::mvrnorm(n=N, mu = m1, Sigma = diag(P)),
                MASS::mvrnorm(n=N, mu = m2, Sigma = diag(P)),
                MASS::mvrnorm(n=N, mu = m3, Sigma = diag(P)))
Xval <- rbind(MASS::mvrnorm(n=N, mu = m1, Sigma = diag(P)),
              MASS::mvrnorm(n=N, mu = m2, Sigma = diag(P)),
              MASS::mvrnorm(n=N, mu = m3, Sigma = diag(P)))

# Generate the labels
Ytrain <- rep(1:3,each=N)
Yval <- rep(1:3,each=N)

# Train the classifier and increase the sparsity parameter from the default # so we penalize more for non-sparse solutions.
res <- accSDA::SZVDcv(cbind(Ytrain,Xtrain),cbind(Yval,Xval),num_gammas=4,
g_mults = c(0,1),beta=2.5,
D=diag(P), maxits=100,tol=list(abs=1e-3,rel=1e-3), k = 3,
ztol=1e-4,sparsity_pen=0.3,quiet=FALSE,penalty=TRUE,scaling=TRUE)
\end{verbatim}

Description

Applies alternating direction methods of multipliers to solve sparse zero variance discriminant analysis.

Usage

\texttt{SZVD\_kFold\_cv(X, \ldots)}

## Default S3 method:
\texttt{SZVD\_kFold\_cv(X, Y, folds, gamms, beta, D, q, maxits, tol,}
\texttt{ztol, feat, penalty, quiet)}
Arguments

- **X**
  - n by p data matrix, variables should be scaled to by sd

- **...**
  - Parameters passed to SZVD.default.

- **Y**
  - n by K indicator matrix.

- **folds**
  - number of folds to use in K-fold cross-validation.

- **gams**
  - Number of regularly spaced regularization parameters to try in [0,1]*max_gamma.
  - See details for how max_gamma is computed in the function.

- **beta**
  - Augmented Lagrangian parameter. Must be greater than zero.

- **D**
  - Penalty dictionary basis matrix.

- **q**
  - Desired number of discriminant vectors.

- **maxits**
  - Number of iterations to run ADMM algorithm.

- **tol**
  - Stopping tolerances for ADMM, must have tol$rel and tol$sabs.

- **ztol**
  - Rounding tolerance for truncating entries to 0.

- **feat**
  - Maximum fraction of nonzero features desired in validation scheme.

- **penalty**
  - Controls whether to apply reweighting of l1-penalty (using sigma = within-class std devs)

- **quiet**
  - toggles between displaying intermediate statistics.

Details

- Add how maxGamma is calculated from the ZVD solution. This function might require a wrapper similar to ASDA.

Value

- SZVDcv returns an object of class "SZVDcv" including a list with the named components DVs and gambest. Where DVs are the discriminant vectors for the best l1 regularization parameter and gambest is the best regularization parameter found in the cross-validation.

- NULL

See Also

- Used by: SZVDcv.

Examples

```r
P <- 150 # Number of variables
N <- 20 # Number of samples per class

# Mean for classes, they are zero everywhere except the first 3 coordinates
m1 <- rep(0,P)
m1[1] <- 3

m2 <- rep(0,P)
m2[2] <- 3
```
m3 <- rep(0, P)
m3[3] <- 3

# Sample dummy data
Xtrain <- rbind(MASS::mvrnorm(n=N, mu = m1, Sigma = diag(P)),
                MASS::mvrnorm(n=N, mu = m2, Sigma = diag(P)),
                MASS::mvrnorm(n=N, mu = m3, Sigma = diag(P)))

# Generate the labels
Ytrain <- cbind(c(rep(1, N), rep(0, 2*N)),
                c(rep(0, N), rep(1, N), rep(0, N)),
                c(rep(0, 2*N), rep(1, N)))

# Normalize the data
Xt <- accSDA::normalize(Xtrain)
Xtrain <- Xt$Xc

# Train the classifier and increase the sparsity parameter from the default
# so we penalize more for non-sparse solutions.
res <- accSDA::SZVD_kFold_cv(Xtrain, Ytrain, folds=2, gams=2, beta=2.5, q=1, D=diag(P),
                             maxits=50, tol=list(abs=1e-2, rel=1e-2),
                             ztol=1e-3, feat=0.3, quiet=FALSE, penalty=TRUE)

---

ZVD

Zero Variance Discriminant Analysis

Description

Implements the ZVD algorithm to solve discriminant vectors.

Usage

ZVD(A, ...)

## Default S3 method:
ZVD(A, scaling = FALSE, get_DVs = FALSE)

Arguments

A                       Matrix, where first column corresponds to class labels.
...                      Parameters passed to ZVD.default.
scaling                 Logical whether to rescale data so each feature has variance 1.
get_DVs                 Logical whether to obtain unpenalized zero-variance discriminant vectors.

Details

This function should potentially be made internal for the release.
Value

SZVDcv returns an object of class "ZVD" including a list with the following named components:

dvs  discriminant vectors (optional).
b  sample between-class covariance.
w  sample within-class covariance.
n  basis for the null space of the sample within-class covariance.
mu  training mean and variance scaling/centering terms
means  vectors of sample class-means.
k  number of classes in given data set.
labels  list of classes.
obs  matrix of data observations.
class_obs  Matrices of observations of each class.

NULL

See Also

Used by: SZVDcv.

Examples

# Generate Gaussian data on three classes with bunch of redundant variables

P <- 300  # Number of variables
N <- 50   # Number of samples per class

# Mean for classes, they are zero everywhere except the first 3 coordinates
m1 <- rep(0,P)
m1[1] <- 3

m2 <- rep(0,P)
m2[2] <- 3

m3 <- rep(0,P)
m3[3] <- 3

# Sample dummy data
Xtrain <- rbind(MASS::mvrnorm(n=N,mu = m1, Sigma = diag(P)),
            MASS::mvrnorm(n=N,mu = m2, Sigma = diag(P)),
            MASS::mvrnorm(n=N,mu = m3, Sigma = diag(P)))

# Generate the labels
Ytrain <- rep(1:3,each=N)

# Normalize the data
Xt <- accSDA::normalize(Xtrain)
Xtrain <- Xt$Xc

# Train the classifier and increase the sparsity parameter from the default
# so we penalize more for non-sparse solutions.
res <- accSDA::ZVD(cbind(Ytrain,Xtrain))
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