

Package ‘regtools’

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Title Regression Tools

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Depends FNN,mvtnorm,dummies,car

Suggests knitr, rmarkdown

VignetteBuilder knitr

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Description Tools for linear, nonlinear and nonparametric regression and classification. Parametric fit assessment using nonparametric methods. One vs. All and All vs. All multiclass classification. Nonparametric regression for general dimension, locally-linear option. Nonlinear regression with Eickert-White method for dealing with heteroscedasticity, k-NN for general dimension and general descriptive functions.

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avalogtrn,avalogpred,ovalogtrn,ovalogpred,knntrn,predict.ovaknn
Classification with More Than 2 Classes

Description

One vs. All, All vs. All tools for multiclass classification, parametric and nonparametric.

Usage

```
ovalogtrn(m, trnxy, truepriors=NULL)
ovalogpred(coefmat, predx)
avalogtrn(m, trnxy)
avalogpred(m, coefmat, predx)
knntrn(y, xdata, m, k, truepriors=NULL)
## S3 method for class 'ovaknn'
predict(object, ...)
classadjjust(econdprobs, wrongratio, trueratio)
```

Arguments

trnxy	Data matrix, one data point per row, Y in the last column.
object	Needed for consistency with generic.
...	Needed for consistency with generic.
y	Vector of response variable data in the training set, with codes 0,1,...m-1.
xdata	X and associated neighbor indices. Output of preprocessx.
coefmat	Output from ovalogtrn.
k	Number of nearest neighbors.
predx	One data point to be predicted.
m	Number of classes in multiclass setting.
econdprobs	Estimated conditional class probabilities, given the predictors.
wrongratio	Incorrect, data-provenanced, $p/(1-p)$, with p being the unconditional probability of a certain class.
trueratio	Same as wrongratio, but with the correct value.
truepriors	True unconditional class probabilities, typically obtained externally.

Details

These functions do classification in the multiclass setting, using the One vs. All method. In addition to logit, the k-Nearest Neighbor method is available. For this, preprocessx must first be called. In the logit case, All vs. All is also offered.

The functions ovalogtrn, avalogtrn and knntrn are used on the training set, and then fed into the prediction functions, ovalogpred, avalogpred and predict.ovaknn. The arguments for the latter are the output of knntrn and a matrix of prediction points (internally referred to as predpts in the code), one per row.

Value

The prediction functions, *ovalogpred*, *avalogpred* and *predict.ovaknn*, return the predicted classes for the points in *predx* or *predpts*.

The functions *ovalogtrn* and *avalogtrn* return the estimated logit coefficient vectors, one per column. There are *m* of them in the former case, $m-1/2$ in the latter, in which case the order of the R function *combin* is used.

The function *knntrn* returns a copy of the *xdata* input, but with an extra component added. The latter is the matrix of estimated regression function values; the element in row *i*, column *j*, is the probability that $Y = j$ given that $X = \text{row } i$ in the *X* data.

Author(s)

Norm Matloff

Examples

```
## Not run:
# toy example, kNN
set.seed <- 9999
x <- runif(25)
y <- sample(0:2,25,replace=TRUE)
xd <- preprocessx(x,2)
kout <- knntrn(y,xd,m=3,k=2)
kout$regist # row 1: 0.5,0.5,0.0
predict(kout,matrix(c(0.25,0.55,0.88),ncol=1)) # 2,0,0

# sim data, kNN
set.seed <- 9999
n <- 1500
# within-grp cov matrix
cv <- rbind(c(1,0.2),c(0.2,1))
xy <- NULL
for (i in 1:3)
  xy <- rbind(xy,rmvnorm(n,mean=rep(i*2.0,2),sigma=cv))
y <- rep(0:2,each=n)
xy <- cbind(xy,y)
xdata <- preprocessx(xy[,-3],20)
oo <- knntrn(y,xdata,m=3,k=20)
predout <- predict(oo,xy[,-3])
mean(predout$predy == y) # about 0.87

\dontrun{
library(mlbench)
data(Vehicle)
xdata <- preprocessx(Vehicle[,-19],25)
kout <- knntrn(Vehicle$Class,xdata,k=25)
predict(kout,as.matrix(Vehicle[1,-19])) # predicted Y is 3

# UCI Letter Recognition data
data(LetterRecognition)
```

```

# prep data
lr <- LetterRecognition
# code Y values
lr[,1] <- as.numeric(lr[,1]) - 1
# training and test sets
lrtrn <- lr[1:14000,]
lrtest <- lr[14001:20000,]
# kNN
xdata <- preprocessx(lrtrn[, -1], 50)
# without setting priors
trnout <- knntrn(lrtrn[, 1], xdata, 26, 50)
ypred <- predict(trnout, as.matrix(lrtest[, -1]))
# how well did it work?
mean(ypred$predy == lrtest[, 1]) # 0.86
# logit
ologout <- ovalogtrn(26, lr[, c(2:17, 1)])
ypred <- ovalogpred(ologout, lr[, -1])
mean(ypred == lr[, 1]) # only 0.73
# try quadratic terms
for (i in 2:17)
  lr <- cbind(lr, lr[, i]^2)
ologout1 <- ovalogtrn(26, lr[, c(2:33, 1)])
ypred <- ovalogpred(ologout1, lr[, -1])
mean(ypred == lr[, 1]) # increased to 0.81
}

## End(Not run)

```

knnest, meany, vary, loclin, predict.knn, preprocessx, kmin, parvsnonparplot, nonparvsxplot, 11, 12
Nonparametric Regression and Classification

Description

Full set of tools for k-NN regression and classification, including both for direct usage and as tools for assessing the fit of parametric models.

Usage

```

knnest(y, xdata, k, nearf=meany)
preprocessx(x, kmax, xval=FALSE)
meany(predpt, nearxy)
vary(predpt, nearxy)
loclin(predpt, nearxy)
## S3 method for class 'knn'
predict(object, ...)
kmin(y, xdata, lossftn=12, nk=5, nearf=meany)

```

```
parvsnonparplot(lmout, knnout, cex=1.0)
nonparvsxplot(knnout, lmout=NULL)
nonparvarplot(knnout)
l2(y, muhat)
l1(y, muhat)
```

Arguments

<code>y</code>	Response variable data in the training set. Vector or matrix, the latter case for vector-valued response, e.g. multiclass classification.
<code>x</code>	X data, predictors, one row per data point, in the training set.
<code>...</code>	Needed for consistency with generic. See Details below for ‘arguments.’
<code>xdata</code>	X and associated neighbor indices. Output of <code>preprocessx</code> .
<code>k</code>	Number of nearest neighbors
<code>object</code>	Output of <code>knnest</code> .
<code>predpt</code>	One point on which to predict, as a vector.
<code>nearxy</code>	A set of X neighbors of a point.
<code>nearf</code>	Function to apply to the nearest neighbors of a point.
<code>kmax</code>	Maximal number of nearest neighbors to find.
<code>xval</code>	Cross-validation flag. If TRUE, then the set of nearest neighbors of a point will not include the point itself.
<code>lossftn</code>	Loss function to be used in cross-validation determination of "best" k.
<code>nk</code>	Number of values of k to try in cross-validation.
<code>lmout</code>	Output of <code>lm</code> .
<code>knnout</code>	Output of <code>knnest</code> .
<code>cex</code>	R parameter to control dot size in plot.
<code>muhat</code>	Vector of estimated regression function values.

Details

The `knnest` function does k-nearest neighbor regression function estimation, in any dimension, i.e. any number of predictor variables, and any number of response variables. This of course includes classification problems case; a scalar $Y = 0,1$ would represent two classes, with the regression function reducing to the conditional probability of class 1, given the predictors.

The `preprocessx` function does the prep work. For each row in `x`, the code finds the `kmax` closest rows to that row. By separating this computation from `knnest`, one can save a lot of overall computing time. If for instance one wants to try the number of nearest neighbors `k` at 25, 50 and 100, one can call `preprocessx` with `kmax` equal to 100, then reuse the results; in calling `knnest` for several values of `k`, we do not need to call `preprocessx` again.

In addition to averaging the neighbor `Y` values, one can specify other types of smoothing by proper specification of the `nearf` function. For instance, one specifies local linear smoothing by setting `nearf` to `loclin`, one could check heteroscedasticity in linear regression models by setting `nearf` to `vary` to estimate conditional variance.

The X , i.e. predictor, data will be scaled by the code, so as to put all predictor variables on an equal footing. The scaling parameters will be recorded, and then applied later in prediction.

One can choose the number of nearest neighbors k through the `kmin` function.

The function `predict.knn` uses the output of `knnest` to do estimation or prediction on new points. Since the output of `knnest` is of class "knn", one invokes this function with the simpler `predict`. The second argument is the set of new points, supplied as a matrix, called `predpts` within the code. A "1-NN" method is used here: Given a new point u whose "Y" value we wish to predict, the code finds the single closest row in the training set, and returns the previously-estimated regression function value at that row.

The functions `ovaknntrn` and `ovaknnpred` are multiclass wrappers for `knnest` and `knnpred`. Here y is coded $0, 1, \dots, m-1$ for the m classes.

The tools here can be useful for fit assessment of parametric models. The `parvsnonparplot` function plots fitted values of parametric model vs. k-NN fitted, `nonparvsxplot` k-NN fitted values against each predictor, one by one.

The functions `l2` and `l1` are used to define L2 and L1 loss.

Value

The return value of `preprocessx` is an R list. Its `x` component is the scaled x matrix, with the scaling factors being recorded in the `scaling` component. The `idxs` component contains the indices of the nearest neighbors of each point in the predictor data, stored in a matrix with `nrow(x)` rows and `k` columns. Row i contains the indices of the nearest rows in x to row i of x . The first of these indices is for the closest point, then for the second-closest, and so on. If cross-validation is requested (`xval = TRUE`, then any point will not be considered a neighbor of itself.

The `knnest` function returns an expanded version of `xdata`, with the expansion consisting of a new component `regest`, the estimated regression function values at the training set points.

The function `predict.knn` returns the predicted Y values at `predpts`. It is called simply via `predict`.

The function `kmin` returns an R list, with the component `meanerrs` containing the cross-validated mean loss function values and `ks` containing the corresponding values of k ; `plot.knn` then plots the former against the latter.

Author(s)

Norm Matloff

Examples

```
set.seed(9999)
x <- matrix(sample(1:100,30),ncol=3)
xd <- preprocessx(x[,1],2,TRUE) # just 1 predictor
ko <- knnest(x[,2],xd,2) # Y is x[,2]
ko$regest # 1st element = 74.5
predict(ko,matrix(76)) # 47.5
ko <- knnest(x[,-1],xd,2) # Y bivar
ko$regest # 1st row = (74.5,31.5)
predict(ko,matrix(76)) # 47.5, 65.0
```

```

xe <- matrix(rnorm(30000), ncol=3)
xe[,-3] <- xe[,-3] + 2
# xe is 2 predictors and epsilon
y <- xe %*% c(1, 0.5, 0.2) # Y
x <- xe[,-3] # X
xdata <- preprocessx(x, 500) # k as high as 500
zout <- knnest(y, xdata, 200)
predict(zout, matrix(c(1, 1), nrow=1)) # about 1.55

## Not run:
data(prgeng)
pe <- prgeng
# dummies for MS, PhD
pe$ms <- as.integer(pe$educ == 14)
pe$phd <- as.integer(pe$educ == 16)
# computer occupations only
pecs <- pe[pe$occ >= 100 & pe$occ <= 109,]
# for simplicity, let's choose a few predictors
pecs1 <- pecs[, c(1, 7, 9, 12, 13, 8)]
# will predict wage income from age, gender etc.
# prepare nearest-neighbor data, k up to 50
xdata <- preprocessx(pecs1[, 1:5], 50)
zout <- knnest(pecs1[, 6], xdata, 50) # k = 50
# find the est. mean income for 42-year-old women, 52 weeks worked, with
# a Master's
predict(zout, matrix(c(42, 2, 52, 0, 0), nrow=1)) # 62106
# try k = 25; don't need to call preprocessx() again
zout <- knnest(pecs1[, 6], xdata, 25)
predict(zout, matrix(c(42, 2, 52, 0, 0), nrow=1)) # 69104
# quite a difference; what k values are good?
kmin(pecs1[, 6], xdata) # at least 50
# what about a man?
zout <- knnest(pecs1[, 6], xdata, 50)
predict(zout, matrix(c(42, 1, 52, 0, 0), nrow=1)) # 78588
# form training and test sets, fit on the former and predict on the
# latter
fullidxs <- 1:nrow(pecs1)
train <- sample(fullidxs, 10000)
xdata <- preprocessx(pecs1[train, 1:5], 50)
trainout <- knnest(pecs1[train, 6], xdata, 50)
testout <- predict(trainout, as.matrix(pecs1[-train, -6]))
# find mean abs. prediction error (about $25K)
mean(abs(pecs1[-train, 6] - testout))
# examples of fit assessment
# look for nonlinear relations between Y and each X
nonparvsxplot(zout) # keep hitting Enter for next plot
# there seem to be quadratic relations with age and wkswrkd, so add quad
# terms and run lm()
pecs2 <- pecs1
pecs2$age2 <- pecs1$age^2
pecs2$wks2 <- pecs1$wkswrkd^2
lmout2 <- lm(wageinc ~ ., data=pecs2)

```

```

# check parametric fit by comparing to kNN
parvsnonparplot(lmout2,zout)
# linear model line somewhat faint, due to large n;
# parametric model seems to overpredict at high end;
# to deal with faintness, reduce size of points
parvsnonparplot(lmout2,zout,cex=0.1)
# assess homogeneity of conditional variance
nonparvarplot(zout)
# hockey stick!

## End(Not run)

# Y vector-valued (3 classes)
# 3 clusters, equal wts, coded 0,1,2
n <- 1500
# within-grp cov matrix
cv <- rbind(c(1,0.2),c(0.2,1))
xy <- NULL
for (i in 1:3)
  xy <- rbind(xy,rmvnorm(n,mean=rep(i*2.0,2),sigma=cv))
y <- rep(0:2,each=n)
xy <- cbind(xy,dummy(y))
xdata <- preprocessx(xy[,-(3:5)],20) # X is xy[,1:2], k <= 20
ko <- knnest(xy[,3:5],xdata,20)
# find predicted Y for each data pt
mx <- apply(as.matrix(ko$regest),1,which.max) - 1
# overall correct classification rate
mean(mx == y) # should be about 0.87

```

```

lmac,makeNA,coef.lmac,vcov.lmac,pcac,loglinac,tbltofakedf
Available Cases Method for Missing Data

```

Description

Various estimators that handle missing data via the Available Cases Method

Usage

```

lmac(xy,nboot=0)
makeNA(m,probna)
## S3 method for class 'lmac'
coef(object,...)
## S3 method for class 'lmac'
vcov(object,...)
pcac(indata,scale=FALSE)
loglinac(x,margin)
tbltofakedf(tbl)

```


Arguments

<code>xy</code>	Matrix or data frame, X values in the first columns, Y in the last column.
<code>indata</code>	Matrix or data frame.
<code>x</code>	Matrix or data frame, one column per variable.
<code>nboot</code>	If positive, number of bootstrap samples to take.
<code>probna</code>	Probability that an element will be NA.
<code>scale</code>	If TRUE, call <code>cor</code> instead of <code>cov</code> .
<code>tbl</code>	Matrix or data frame, one column per variable.
<code>tbl</code>	An R table.
<code>m</code>	Number of synthetic NAs to insert.
<code>object</code>	Output from <code>lmac</code> .
<code>...</code>	Needed for consistency with generic function. Not used.
<code>margin</code>	A list of vectors specifying the model, as in <code>loglin</code> .

Details

The Available Cases (AC) approach applies to statistical methods that depend only on products of k of the variables, so that cases having non-NA values for those k variables can be used, as opposed to using only cases that are fully intact in all variables, the Complete Cases (CC) approach. In the case of linear regression, for instance, the estimated coefficients depend only on covariances between the variables (both predictors and response). This approach assumes that the cases with missing values have the same distribution as the intact cases.

The `lmac` function forms OLS estimates as with `lm`, but applying AC, in contrast to `lm`, which uses the CC method.

The `pcac` function is an AC substitute for `prcomp`. The data is centered, corresponding to a fixed value of `center = TRUE` in `prcomp`. It is also scaled if `scale = TRUE`, corresponding to `scale. = TRUE` in `prcomp`; . Due to AC, there is a small chance of negative eigenvalues, in which case `stop` will be called.

The `loglinac` function is an AC substitute for `loglin`. The latter takes tables as input, but `loglinac` takes the raw data. If you have just the table, use `tbltofakedf` to regenerate a usable data frame.

The `makeNA` function is used to insert random NA values into data, for testing purposes.

Value

For `lmac`, an object of class `'lmac'`, with components

- coefficients, as with `lm`; accessible directly or by calling `coef`, as with `lm`
- `fitted.values`, as with `lm`
- residuals, as with `lm`
- `r2`, (unadjusted) R-squared
- `cov`, for `nboot > 0` the estimated covariance matrix of the vector of estimated regression coefficients; accessible directly or by calling `vcov`, as with `lm`

For `pcac`, an R list, with components

- `sdev`, as with `prcomp`
- `rotation`, as with `prcomp`

For `loglinac`, an R list, with components

- `param`, estimated coefficients, as in `loglin`
- `fit`, estimated expected call counts, as in `loglin`

Author(s)

Norm Matloff

Examples

```
n <- 25000
w <- matrix(rnorm(2*n),ncol=2) # x and epsilon
x <- w[,1]
y <- x + w[,2]
# insert some missing values
nmiss <- round(0.1*n)
x[sample(1:n,nmiss)] <- NA
nmiss <- round(0.2*n)
y[sample(1:n,nmiss)] <- NA
acout <- lmac(cbind(x,y))
coef(acout) # should be near pop. values 0 and 1
```

ltrfreqs

Letter Frequencies

Description

This is data consists of capital letter frequencies obtained at <http://www.math.cornell.edu/~mec/2003-2004/cryptography/subs/frequencies.html>

Usage

```
data(ltrfreqs); ltrfreqs
```

 mm *Method of Moments, Including Possible Regression Terms*

Description

Method of Moments computation for almost any statistical problem that has derivatives with respect to theta. Capable of handling models that include parametric regression terms, but not need be a regression problem. (This is not *Generalized* Method of Moments; see the package **gmm** for the latter.)

Usage

```
mm(m,g,x,init=rep(0.5,length(m)),eps=0.0001,maxiters=1000)
```

Arguments

m	Vector of sample moments, "left-hand sides" of moment equations.
g	Function of parameter estimates, forming the "right-hand sides." This is a multivariate-valued function, of dimensionality equal to that of m.
init	Vector of initial guesses for parameter estimates. If components are named, these will be used as labels in the output.
eps	Convergence criterion.
maxiters	Maximum number of iterations.
x	Input data.

Details

Standard Newton-Raphson methods are used to solve for the parameter estimates, with `numericDeriv` being used to find the approximate derivatives.

Value

R list consisting of components `tht`, the vector of parameter estimates, and `numiters`, the number of iterations performed.

Author(s)

Norm Matloff

Examples

```
x <- rgamma(1000,2)
m <- c(mean(x),var(x))
g <- function(x,theta) { # from theoretical properties of gamma distr.
  g1 <- theta[1] / theta[2]
  g2 <- theta[1] / theta[2]^2
  c(g1,g2)
}
```

```

}
# should output about 2 and 1
mm(m,g,x)

## Not run:
library(mfp)
data(bodyfat)
# model as a beta distribution
g <- function(x,theta) {
  t1 <- theta[1]
  t2 <- theta[2]
  t12 <- t1 + t2
  meanb <- t1 / t12
  m1 <- meanb
  m2 <- t1*t2 / (t12^2 * (t12+1))
  c(m1,m2)
}
x <- bodyfat$brozek/100
m <- c(mean(x),var(x))
# about 4.65 and 19.89
mm(m,g,x)

## End(Not run)

```

nlshc

Heteroscedastic Nonlinear Regression

Description

Extension of nls to the heteroscedastic case.

Usage

```
nlshc(nlsout, type='hc3')
```

Arguments

nlsout	Object of type 'nls'.
type	Eickert-White algorithm to use. See documentation for nls .

Details

Calls nls but then forms a different estimated covariance matrix for the estimated regression coefficients, applying the Eickert-White technique to handle heteroscedasticity. This then gives valid statistical inference in that setting.

Some users may prefer to use nlsLM of the package **minpack.lm** instead of nls. This is fine, as both functions return objects of class 'nls'.

Value

Estimated covariance matrix

Author(s)

Norm Matloff

References

Zeileis A (2006), Object-Oriented Computation of Sandwich Estimators. *Journal of Statistical Software*, **16**(9), 1–16, <http://www.jstatsoft.org/v16/i09/>.

Examples

```
# simulate data from a setting in which mean Y is
# 1 / (b1 * X1 + b2 * X2)
n <- 250
b <- 1:2
x <- matrix(rexp(2*n),ncol=2)
meany <- 1 / (x %>% b) # reg ftn
y <- meany + (runif(n) - 0.5) * meany # heterosced epsilon
xy <- cbind(x,y)
xy <- data.frame(xy)
# see nls() docs
nlout <- nls(X3 ~ 1 / (b1*X1+b2*X2),
  data=xy,start=list(b1 = 1,b2=1))
nlshc(nlout)
```

prgeng

Silicon Valley programmers and engineers

Description

This data set is adapted from the 2000 Census (5% sample, person records). It is restricted to programmers and engineers in the Silicon Valley area.

The variable codes, e.g. occupational codes, are available from the Census Bureau, at <http://www.census.gov/prod/cen2000/doc/pums.pdf>. (Short code lists are given in the record layout, but longer ones are in the appendix Code Lists.)

The variables are:

- age, with a U(0,1) variate added for jitter
- cit, citizenship; 1-4 code various categories of citizens; 5 means noncitizen (including permanent residents)
- educ: 01-09 code no college; 10-12 means some college; 13 is a bachelor's degree, 14 a master's, 15 a professional deal and 16 is a doctorate
- occ, occupation

- birth, place of birth
- wageinc, wage income
- wkswrkd, number of weeks worked
- yrentry, year of entry to the U.S. (0 for natives)
- powpuma, location of work
- gender, 1 for male, 2 for female

Usage

```
data(prgeng); prgeng
```

```
ridgelm,plot.rlm      Ridge Regression
```

Description

Similar to `lm.ridge` in MASS packaged included with R, but with a different kind of scaling and a little nicer plotting.

Usage

```
ridgelm(xy,lambda = seq(0.01,1,0.01),mapback=TRUE)
## S3 method for class 'rlm'
plot(x,y,...)
```

Arguments

<code>xy</code>	Data, response variable in the last column.
<code>lambda</code>	Vector of desired values for the ridge parameter.
<code>mapback</code>	If TRUE, the scaling that had been applied to the original data will be map back to the original scale, so that the estimated regression coefficients are now on the scale of the original data.
<code>x</code>	Object of type 'rlm', output of <code>ridgelm</code> .
<code>y</code>	Needed for consistency with the generic. Not used.
<code>...</code>	Needed for consistency with the generic. Not used.

Details

Centers and scales the predictors X, and centers the response variable Y. Computes $X'X$ and then solves $[(X'X)/n + \lambda I]b = X'Y/n$ for b. The $1/n$ factors are important, making the diagonal elements of $(X'X)/n$ all 1s and thus facilitating choices for the lambdas in a manner independent of the data.

Calling `plot` on the output of `ridgelm` dispatches to `plot.rlm`, thus displaying the ridge traces.

Value

The function `ridgelm` returns an object of class `'rlm'`, with components `bhats`, the estimated beta vectors, one column per lambda value, and `lambda`, a copy of the input.

Author(s)

Norm Matloff

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