

Package ‘APPEstimation’

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Type Package

Title Adjusted Prediction Model Performance Estimation

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Description Calculating predictive model performance measures adjusted for predictor distributions using density ratio method (Sugiyama et al., (2012, ISBN:9781139035613)). L1 and L2 error for continuous outcome and C-statistics for binomial outcome are computed.

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APPEstimation-package *R function to calculate model performance measure adjusted for predictor distributions.*

Description

This package provides the function to estimate model performance measures (L_1 , L_2 , C -statistics). The difference in the distribution of predictors between two datasets (training and validation) is adjusted by a density ratio estimate.

Details

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References

Sugiyama, M., Suzuki, T. & Kanamori, T. Density Ratio Estimation in Machine Learning. Cambridge University Press 2012. ISBN:9781139035613.

Examples

```

set.seed(100)

# generating learning data
n0 = 100
Z = cbind(rbeta(n0, 5, 5), rbeta(n0, 5, 5))
Y = apply(Z, 1, function (xx) {
  rbinom(1, 1, (1/(1+exp(-(sum(c(-2,2,2) * c(1,xx))))))))))
dat = data.frame(Y=Y, Za=Z[,1], Zb=Z[,2])

# the model to be evaluated
mdl = glm(Y~., binomial, data=dat)

# validation dataset, with different centers on predictors
n1 = 100
Z1 = cbind(rbeta(n1, 6, 4), rbeta(n1, 6, 4))
Y1 = apply(Z1, 1, function (xx) {
  rbinom(1, 1, (1/(1+exp(-(sum(c(-2,2,2) * c(1,xx))))))))))
dat1 = data.frame(Y=Y1, Za=Z1[,1], Zb=Z1[,2])

# calculation of L1 and L2 for this model
appe.glm(mdl, dat, dat1, reps=0)

```

appe.glm

C-statistics adjusted for predictor distributions

Description

Calculates adjusted *C* statistics by predictor distributions for a generalized linear model with binary outcome.

Usage

```

appe.glm(mdl, dat.train, dat.test, method = "uLSIF", sigma = NULL,
         lambda = NULL, kernel_num = NULL, fold = 5, stabilize = TRUE,
         qstb = 0.025, reps = 2000, conf.level = 0.95)

```

Arguments

- mdl a glm object describing a prediction model to be evaluated.
- dat.train a dataframe used to construct a prediction model (specified in mdl), corresponding to a training data. Need to include outcome and all predictors.
- dat.test a dataframe corresponding to a validation (testing) data. Need to include outcome and all predictors.
- method uLSIF or KLIEP. Same as the argument in densratio function from densratio package.

<code>sigma</code>	a positive numeric vector corresponding to candidate values of a bandwidth for Gaussian kernel. Same as the argument in <code>densratio</code> function from <code>densratio</code> package.
<code>lambda</code>	a positive numeric vector corresponding to candidate values of a regularization parameter. Same as the argument in <code>densratio</code> function from <code>densratio</code> package.
<code>kernel_num</code>	a positive integer corresponding to number of kernels. Same as the argument in <code>densratio</code> function from <code>densratio</code> package.
<code>fold</code>	a positive integer corresponding to a number of the folds of cross-validation in the KLIEP method. Same as the argument in <code>densratio</code> function from <code>densratio</code> package.
<code>stabilize</code>	a logical value as to whether tail weight stabilization is performed or not. If TRUE, both tails of the estimated density ratio distribution are replaced by the constant value which is specified at <code>qstb</code> option.
<code>qstb</code>	a positive numerical value less than 1 to control the degree of weight stabilization. Default value is 0.025, indicating estimated density ratio values less than the 2.5 percentile and more than the 97.5 percentile are set to 2.5 percentile and 97.5 percentile, respectively.
<code>reps</code>	a positive integer to specify bootstrap repetitions. If 0, bootstrap calculations are not performed.
<code>conf.level</code>	a numerical value indicating a confidence level of interval.

Value

Adjusted and non-adjusted estimates of C -statistics are provided as matrix form. "Cstat" indicates non-adjusted version, "C adjusted by score" indicates adjusted version by linear predictors distribution, and "C adjusted by predictors" indicates adjusted version by predictor distributions (multi-dimensionally). For confidence intervals, "Percentile" indicates a confidence interval by percentile method and "Approx" indicates approximated versions by Normal distribution.

Examples

```
set.seed(100)

# generating learning data
n0 = 100
Z = cbind(rbeta(n0, 5, 5), rbeta(n0, 5, 5))
Y = apply(Z, 1, function (xx) {
  rbinom(1, 1, (1/(1+exp(-(sum(c(-2,2,2) * c(1,xx))))))))))
dat = data.frame(Y=Y, Za=Z[,1], Zb=Z[,2])

# the model to be evaluated
mdl = glm(Y~., binomial, data=dat)

# validation dataset, with different centers on predictors
n1 = 100
Z1 = cbind(rbeta(n1, 6, 4), rbeta(n1, 6, 4))
Y1 = apply(Z1, 1, function (xx) {
```

```

      rbinom(1, 1, (1/(1+exp(-(sum(c(-2,2,2) * c(1,xx))))))))))
dat1 = data.frame(Y=Y1, Za=Z1[,1], Zb=Z1[,2])

# calculation of L1 and L2 for this model
appe.glm mdl, dat, dat1, reps=0)

```

appe.lm

L₁ and L₂ errors adjusted for predictor distributions

Description

Calculates adjusted L_1 and L_2 errors by predictor distributions for a linear model.

Usage

```

appe.lm(mdl, dat.train, dat.test, method = "uLSIF", sigma = NULL,
        lambda = NULL, kernel_num = NULL, fold = 5, stabilize = TRUE,
        qstb = 0.025, reps = 2000, conf.level = 0.95)

```

Arguments

mdl	a lm object describing a prediction model to be evaluated.
dat.train	same as in appe.glm.
dat.test	same as in appe.glm.
method	same as in appe.glm.
sigma	same as in appe.glm.
lambda	same as in appe.glm.
kernel_num	same as in appe.glm.
fold	same as in appe.glm.
stabilize	same as in appe.glm.
qstb	same as in appe.glm.
reps	same as in appe.glm.
conf.level	same as in appe.glm.

Value

Adjusted and non-adjusted estimates of L_1 and L_2 errors are provided as matrix form. "L1" and "L2" indicate non-adjusted versions, "L1 adjusted by score" and "L2 adjusted by score" indicate adjusted versions by linear predictors distribution, "L1 adjusted by predictors" and "L2 adjusted by predictors" indicate adjusted versions by predictor distributions (multi-dimensionally). For confidence intervals, "Percentile" indicates a confidence interval by percentile method and "Approx" indicates approximated versions by Normal distribution.

Examples

```
set.seed(100)

# generating development data
n0 = 100
Z = cbind(rbeta(n0, 3, 3), rbeta(n0, 3, 3))
Y = apply(Z, 1, function(xx) { rlnorm(1, sum(c(1, 1) * xx), 0.3) })
dat = data.frame(Za=Z[,1], Zb=Z[,2], Y=Y)

# the model to be evaluated
mdl = lm(Y~ Za + Zb, data=dat)

# generating validation dataset
n1 = 100
Z1 = cbind(rbeta(n0, 3.5, 2.5), rbeta(n0, 3.5, 2.5))
Y1 = apply(Z1, 1, function(xx) { rlnorm(1, sum(c(1, 1) * xx), 0.3) })
dat1 = data.frame(Za=Z1[,1], Zb=Z1[,2], Y=Y1)

# calculation of L1 and L2 for this model
ape.lm(mdl, dat, dat1, reps=0)
```

cvalest.bin

Estimation of C-statistics

Description

Calculates *C*-statistics. Individual case weight can be incorporated.

Usage

```
cvalest.bin(Y, scr, wgt = NULL)
```

Arguments

Y	a numerical vector of inary outcome, either 0 or 1.
scr	a numerical vector of continuous variable.
wgt	a numerical vector corresponding to individual weight.

Value

C-statistics is provided.

densratio.appe	<i>A wrapper function</i>
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Description

A wrapper function to use "densratio" function from the densratio package.

Usage

```
densratio.appe(xtrain, xtest, method = "uLSIF", sigma = NULL,  
              lambda = NULL, kernel_num = NULL, fold = 5,  
              stabilize = TRUE, qstb = 0.025)
```

Arguments

xtrain	a dataframe used to construct a prediction model.
xtest	a dataframe corresponding to a validation (testing) data.
method	same as in <code>appe.glm</code> .
sigma	same as in <code>appe.glm</code> .
lambda	same as in <code>appe.glm</code> .
kernel_num	same as in <code>appe.glm</code> .
fold	same as in <code>appe.glm</code> .
stabilize	same as in <code>appe.glm</code> .
qstb	same as in <code>appe.glm</code> .

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