

# Package: customizedTraining (via r-universe)

November 5, 2024

**Type** Package

**Title** Customized Training for Lasso and Elastic-Net Regularized  
Generalized Linear Models

**Version** 1.2

**Date** 2019-01-28

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**Imports** FNN, glmnet

**Description** Customized training is a simple technique for transductive learning, when the test covariates are known at the time of training. The method identifies a subset of the training set to serve as the training set for each of a few identified subsets in the training set. This package implements customized training for the glmnet() and cv.glmnet() functions.

**License** GPL-2

**Repository** <https://saberpowers.r-universe.dev>

**RemoteUrl** <https://github.com/saberpowers/customizedtraining>

**RemoteRef** HEAD

**RemoteSha** 8feae96962cb66b2d396c26327ba59ba2a9e836a

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customizedTraining-package  
*customizedTraining*

---

## Description

Fit a regularized lasso model using customized training

## Details

Customized training is a simple strategy for making predictions on test data when the features of the test data are available at the time of model fitting. The method clusters the data to find training points close to each test point and then fits a GLM elastic net model separately in each training cluster. In this way, customized training is a localized method for transductive learning. In contrast with local regression, however, instead of fitting a separate regression model for each test point, customized training fits only one model for each of a handful of clusters.

Use `customizedGlmnet()` to fit a `glmnet()` model with customized training. Use `cv.customizedGlmnet()` to do the same while choosing the regularization parameter and potentially the number of groups using cross-validation. The `plot()` and `predict()` methods are implemented for both `customizedGlmnet()` and `cv.customizedGlmnet()`.

## Author(s)

Scott Powers, Trevor Hastie, Robert Tibshirani

Maintainer: Scott Powers <sspowers@stanford.edu>

## References

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

## See Also

[glmnet](#)

**Examples**

```

require(glmnet)

# Simulate three groups, each with a different sparse linear model
# producing the response, and fit customized training model (with CV) to the
# synthetic data

# Simulation parameters:
n = m = 300
p = 100
q = 10
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

# Produce the synthetic data:
beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
foldid = sample(rep(1:10, length = nrow(x.train)))

# Fit the customized training model with CV:
fit2 = cv.customizedGlmnet(x.train, y.train, x.test, Gs = 1:3,
  family = "gaussian", foldid = foldid)

# Print the optimal number of groups and value of lambda:
fit2$G.min
fit2$lambda.min

# Print the customized training model fit:
fit2

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit2))^2)

# Plot nonzero coefficients by group:
plot(fit2)

```

**Description**

Fit a regularized lasso model using customized training

**Usage**

```
customizedGlmnet(xTrain, yTrain, xTest, groupid = NULL, G = NULL,
  family = c("gaussian", "binomial", "multinomial"), dendrogram = NULL,
  dendrogramTestIndices = NULL)
```

**Arguments**

xTrain	an n-by-p matrix of training covariates
yTrain	a length-n vector of training responses. Numeric for family = "gaussian". Factor or character for family = "binomial" or family = "multinomial"
xTest	an m-by-p matrix of test covariates
groupid	an optional length-m vector of group memberships for the test set. If specified, customized training subsets are identified using the union of nearest neighbor sets for each test group. Either groupid or G must be specified
G	a positive integer indicating the number of clusters for the joint clustering of the test and training data. Ignored if groupid is specified. Either groupid or G must be specified
family	response type
dendrogram	optional output from hclust on the joint covariate data. Used by cv.customizedGlmnet so that clustering is not computed redundantly
dendrogramTestIndices	optional set of indices (corresponding to dendrogram) held out in cross-validation. Used by cv.customizedGlmnet

**Details**

Identify customized training subsets of the training data through one of two methods: (1) If groupid is specified, grouping the test data, then for each test group find the 10 nearest neighbors of each observation in the group and use the union of these nearest neighbor sets as the customized training set or (2) If G is specified, jointly cluster the test and training data using hierarchical clustering with complete linkage. Within each cluster, the training data are used as the customized training subset for the test data. Once the customized training subsets have been identified, use glmnet to fit an l1-regularized regression model to each.

**Value**

	an object with class customizedGlmnet
call	the call that produced this object
CTset	a list containing the customized training subsets for each test group
fit	a list containing the glmnet fit for each test group
groupid	a length-m vector containing the group memberships of the test data

x	a list containing <code>train</code> (which is the input <code>xTrain</code> ) and <code>test</code> (which is the input <code>xTest</code> ). Specified in function call
y	training response vector (specified in function call)
family	response type (specified in function call)
standard	the fit of <code>glmnet</code> to the entire training set using standard training

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**References**

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

**See Also**

[print.customizedGlmnet](#), [predict.customizedGlmnet](#), [plot.customizedGlmnet](#), [cv.customizedGlmnet](#)

**Examples**

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks
```

```

fit1 = customizedGlmnet(x.train, y.train, x.test, G = 3,
  family = "gaussian")

# Print the customized training model fit:
fit1

# Extract nonzero regression coefficients for each group:
nonzero(fit1, lambda = 10)

# Compute test error using the predict function:
mean((y.test - predict(fit1, lambda = 10))^2)

# Plot nonzero coefficients by group:
plot(fit1, lambda = 10)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = customizedGlmnet(x.train, y.train, x.test, group.id)

# Print the customized training model fit:
fit2

# Extract nonzero regression coefficients for each group:
nonzero(fit2, lambda = 10)

# Compute test error using the predict function:
mean((y.test - predict(fit2, lambda = 10))^2)

# Plot nonzero coefficients by group:
plot(fit2, lambda = 10)

```

---

cv.customizedGlmnet    *cross validation for customizedGlmnet*

---

## Description

Does k-fold cross-validation for customizedGlmnet and returns a values for G and lambda

## Usage

```

cv.customizedGlmnet(xTrain, yTrain, xTest = NULL, groupid = NULL, Gs = NULL,
  dendrogram = NULL, dendrogramCV = NULL, lambda = NULL,
  nfolds = 10, foldid = NULL, keep = FALSE,
  family = c("gaussian", "binomial", "multinomial"), verbose = FALSE)

```

**Arguments**

xTrain	an n-by-p matrix of training covariates
yTrain	a length-n vector of training responses. Numeric for family = "gaussian". Factor or character for family = "binomial" or family = "multinomial"
xTest	an m-by-p matrix of test covariates. May be left NULL, in which case cross validation predictions are made internally on the training set and no test predictions are returned.
groupid	an optional length-m vector of group memberships for the test set. If specified, customized training subsets are identified using the union of nearest neighbor sets for each test group, in which case cross-validation is used only to select the regularization parameter lambda, not the number of clusters G. Either groupid or Gs must be specified
Gs	a vector of positive integers indicating the numbers of clusters over which to perform cross-validation to determine the best number. Ignored if groupid is specified. Either groupid or Gs must be specified
dendrogram	optional output from hclust on the joint covariate data. Useful if method is being used several times to avoid redundancy in calculations
dendrogramCV	optional output from hclust on the training covariate data. Used as joint clustering result for cross-validation. Useful to specify in advance if method is being used several times to avoid redundancy in calculations
lambda	sequence of values to use for the regularization parameter lambda. Recommended to leave as NULL and allow glmnet to choose automatically.
nfolds	number of folds – default is 10. Ignored if foldid is specified
foldid	an optional length-n vector of fold memberships used for cross-validation
keep	Should fitted values on the training set from cross validation be included in output? Default is FALSE.
family	response type
verbose	Should progress be printed to console as folds are evaluated during cross-validation? Default is FALSE.

**Value**

	an object of class cv.customizedGlmnet
call	the call that produced this object
G.min	unless groupid is specified, the number of clusters minimizing CV error
lambda	the sequence of values of the regularization parameter lambda considered
lambda.min	the value of the regularization parameter lambda minimizing CV error
error	a matrix containing the CV error for each G and lambda
fit	a customizedGlmnet object fit using G.min and lambda.min. Only returned if xTest is not NULL.
prediction	a length-m vector of predictions for the test set, using the tuning parameters which minimize cross-validation error. Only returned if xTest is not NULL.

`selected` a list of nonzero variables for each customized training set, using `G.min` and `lambda.min`. Only returned if `xTest` is not `NULL`.

`cv.fit` a array containing fitted values on the training set from cross validation. Only returned if `keep` is `TRUE`.

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**References**

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

**See Also**

[customizedGlmnet](#), [plot.cv.customizedGlmnet](#), [predict.cv.customizedGlmnet](#)

**Examples**

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]
foldid = sample(rep(1:10, length = nrow(x.train)))

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = cv.customizedGlmnet(x.train, y.train, x.test, Gs = c(1, 2, 3, 5),
```



```
    family = "gaussian", foldid = foldid)

# Print the optimal number of groups and value of lambda:
fit1$G.min
fit1$lambda.min

# Print the customized training model fit:
fit1

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit1))^2)

# Plot nonzero coefficients by group:
plot(fit1)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
foldid = apply(z == 1, 1, which)[1:n]
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = cv.customizedGlmnet(x.train, y.train, x.test, group.id, foldid = foldid)

# Print the optimal value of lambda:
fit2$lambda.min

# Print the customized training model fit:
fit2

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit2))^2)

# Plot nonzero coefficients by group:
plot(fit2)

# Example 3: If there is no test set, but the training set is organized into
# blocks, you can do cross validation with these blocks as the basis for the
# customized training sets.

fit3 = cv.customizedGlmnet(x.train, y.train, foldid = foldid)

# Print the optimal value of lambda:
fit3$lambda.min

# Print the customized training model fit:
fit3

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit3))^2)

# Plot nonzero coefficients by group:
plot(fit3)
```

---

nonzero	<i>return selected variables</i>
---------	----------------------------------

---

**Description**

nonzero is a generic function for returning the set of variables selected by a model

**Usage**

```
nonzero(object, ...)
```

**Arguments**

object	a model object for which the set of selected variables is desired
...	additional arguments, e.g. tuning parameters

**Value**

The form of the value returned by nonzero depends on the class of its argument. See the documentation of the particular methods for details of what is produced by that method.

**See Also**

[nonzero.customizedGlmnet](#), [nonzero.singleton](#)

---

nonzero.customizedGlmnet	<i>return selected variables from a customizedGlmnet object</i>
--------------------------	---

---

**Description**

Returns a list of vectors of selected variables for each separate glmnet model fit by a customizedGlmnet object

**Usage**

```
## S3 method for class 'customizedGlmnet'
nonzero(object, lambda = NULL, ...)
```

**Arguments**

object	fitted customizedGlmnet model object
lambda	value of regularization parameter to use. Must be specified
...	ignored

**Value**

a list of vectors, each vectors representing one of the glmnet models fit by the customizedGlmnet model. Each vector gives the indices of the variables selected by the model

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**References**

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

**See Also**

[nonzero](#), [customizedGlmnet](#)

**Examples**

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = customizedGlmnet(x.train, y.train, x.test, G = 3,
  family = "gaussian")
```

```
# Extract nonzero regression coefficients for each group:
nonzero(fit1, lambda = 10)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = customizedGlmnet(x.train, y.train, x.test, group.id)

# Extract nonzero regression coefficients for each group:
nonzero(fit2, lambda = 10)
```

---

nonzero.singleton      *return selected variables from a singleton object*

---

### Description

Returns NULL. Intended for internal use only

### Usage

```
## S3 method for class 'singleton'
nonzero(object, ...)
```

### Arguments

object	an object of class singleton
...	ignored

### Value

NULL

### See Also

[nonzero](#)

---

plot.customizedGlmnet *visualize variables selected in each customized training subset*

---

### Description

Produces a plot, with a row for each customized training submodel, showing the variables selected in the subset, with variables along the horizontal axis

### Usage

```
## S3 method for class 'customizedGlmnet'  
plot(x, lambda, ...)
```

### Arguments

x	a fitted customizedGlmnet object
lambda	regularization parameter. Required
...	ignored

### Author(s)

Scott Powers, Trevor Hastie, Robert Tibshirani

### See Also

[plot, customizedGlmnet](#)

### Examples

```
require(glmnet)  
  
# Simulate synthetic data  
  
n = m = 150  
p = 50  
q = 5  
K = 3  
sigmaC = 10  
sigmaX = sigmaY = 1  
set.seed(5914)  
  
beta = matrix(0, nrow = p, ncol = K)  
for (k in 1:K) beta[sample(1:p, q), k] = 1  
c = matrix(rnorm(K*p, 0, sigmaC), K, p)  
eta = rnorm(K)  
pi = (exp(eta)+1)/sum(exp(eta)+1)  
z = t(rmultinom(m + n, 1, pi))  
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)  
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)
```

```

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = customizedGlmnet(x.train, y.train, x.test, G = 3,
  family = "gaussian")

# Plot nonzero coefficients by group:
plot(fit1, lambda = 10)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = customizedGlmnet(x.train, y.train, x.test, group.id)

# Plot nonzero coefficients by group:
plot(fit2, lambda = 10)

```

---

```
plot.cv.customizedGlmnet
```

*visualize variables selected in each customized training subset, from a cross-validated model*

---

## Description

Produces a plot, with a row for each customized training submodel, showing the variables selected in the subset, with variables along the horizontal axis. The lambda used is the one which minimizes CV error.

## Usage

```
## S3 method for class 'cv.customizedGlmnet'
plot(x, ...)
```

## Arguments

```
x          a fitted cv.customizedGlmnet object
...        ignored
```

## Author(s)

Scott Powers, Trevor Hastie, Robert Tibshirani

**See Also**[plot, cv.customizedGlmnet](#)**Examples**

```

require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]
foldid = sample(rep(1:10, length = nrow(x.train)))

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = cv.customizedGlmnet(x.train, y.train, x.test, Gs = c(1, 2, 3, 5),
  family = "gaussian", foldid = foldid)

# Print the optimal number of groups and value of lambda:
fit1$G.min
fit1$lambda.min

# Print the customized training model fit:
fit1

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit1))^2)

# Plot nonzero coefficients by group:
plot(fit1)

```

```

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
foldid = apply(z == 1, 1, which)[1:n]
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = cv.customizedGlmnet(x.train, y.train, x.test, group.id, foldid = foldid)

# Print the optimal value of lambda:
fit2$lambda.min

# Print the customized training model fit:
fit2

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit2))^2)

# Plot nonzero coefficients by group:
plot(fit2)

# Example 3: If there is no test set, but the training set is organized into
# blocks, you can do cross validation with these blocks as the basis for the
# customized training sets.

fit3 = cv.customizedGlmnet(x.train, y.train, foldid = foldid)

# Print the optimal value of lambda:
fit3$lambda.min

# Print the customized training model fit:
fit3

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit3))^2)

# Plot nonzero coefficients by group:
plot(fit3)

```

---

```
predict.customizedGlmnet
```

*make predictions from a customizedGlmnet object*

---

## Description

Returns predictions for the test set provided at the time of fitting the customizedGlmnet object.

## Usage

```
## S3 method for class 'customizedGlmnet'
```



```
predict(object, lambda,
        type = c('response', 'class'), ...)
```

### Arguments

object	a fitted customizedGlmnet object
lambda	regularization parameter
type	Type of prediction, currently only "response" and "class" are supported. Type "response" returns fitted values for "gaussian" family and fitted probabilities for "binomial" and "multinomial" families. Type "class" applies only to "binomial" and "multinomial" families and returns the class with the highest fitted probability.
...	ignored

### Value

a vector of predictions corresponding to the test data input to the model at the time of fitting

### Author(s)

Scott Powers, Trevor Hastie, Robert Tibshirani

### References

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

### See Also

[predict, customizedGlmnet](#)

### Examples

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
```

```

z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = customizedGlmnet(x.train, y.train, x.test, G = 3,
  family = "gaussian")

# Compute test error using the predict function:
mean((y.test - predict(fit1, lambda = 10))^2)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = customizedGlmnet(x.train, y.train, x.test, group.id)

# Compute test error using the predict function:
mean((y.test - predict(fit2, lambda = 10))^2)

```

---

`predict.cv.customizedGlmnet`

*make predictions from a `cv.customizedGlmnet` object*

---

## Description

Returns predictions for test set provided at time of fitting, using regularization parameter which minimizes CV error

## Usage

```
## S3 method for class 'cv.customizedGlmnet'
predict(object, ...)
```

## Arguments

<code>object</code>	a fitted <code>cv.customizedGlmnet</code> object
<code>...</code>	additional arguments to be passed to <code>predict.customizedGlmnet</code>

**Value**

a vector of predictions corresponding to the test set provided when the model was fit. The results are for the regularization parameter chosen by cross-validation

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**References**

Scott Powers, Trevor Hastie and Robert Tibshirani (2015) "Customized training with an application to mass spectrometric imaging of gastric cancer data." *Annals of Applied Statistics* 9, 4:1709-1725.

**See Also**

[predict, cv.customizedGlmnet](#)

**Examples**

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]
foldid = sample(rep(1:10, length = nrow(x.train)))

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = cv.customizedGlmnet(x.train, y.train, x.test, Gs = c(1, 2, 3, 5),
  family = "gaussian", foldid = foldid)
```

```
# Print the optimal number of groups and value of lambda:
fit1$G.min
fit1$lambda.min

# Print the customized training model fit:
fit1

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit1))^2)

# Plot nonzero coefficients by group:
plot(fit1)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
foldid = apply(z == 1, 1, which)[1:n]
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = cv.customizedGlmnet(x.train, y.train, x.test, group.id, foldid = foldid)

# Print the optimal value of lambda:
fit2$lambda.min

# Print the customized training model fit:
fit2

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit2))^2)

# Plot nonzero coefficients by group:
plot(fit2)

# Example 3: If there is no test set, but the training set is organized into
# blocks, you can do cross validation with these blocks as the basis for the
# customized training sets.

fit3 = cv.customizedGlmnet(x.train, y.train, foldid = foldid)

# Print the optimal value of lambda:
fit3$lambda.min

# Print the customized training model fit:
fit3

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit3))^2)

# Plot nonzero coefficients by group:
plot(fit3)
```

---

predict.singleton      *make predictions from a “singleton” object*

---

**Description**

Returns the value stored in the singleton. Intended for internal use only.

**Usage**

```
## S3 method for class 'singleton'  
predict(object, type = c('response', 'class', 'nonzero'),  
  ...)
```

**Arguments**

object	an object of class singleton
type	Type of prediction to be returned, "response" or "class"
...	ignored

**Value**

The value of the singleton

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**See Also**

[predict](#)

---

print.customizedGlmnet  
*print the summary of a fitted customizedGlmnet object*

---

**Description**

Print the numbers of training observations and test observations in each submodel of the customizedGlmnet fit

**Usage**

```
## S3 method for class 'customizedGlmnet'  
print(x, ...)
```

**Arguments**

x                    fitted customizedGlmnet object  
 ...                  ignored

**Author(s)**

Scott Powers, Trevor Hastie, Robert Tibshirani

**See Also**

[print, customizedGlmnet](#)

**Examples**

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = customizedGlmnet(x.train, y.train, x.test, G = 3,
  family = "gaussian")

# Print the customized training model fit:
fit1
```

```
# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = customizedGlmnet(x.train, y.train, x.test, group.id)

# Print the customized training model fit:
fit2
```

---

```
print.cv.customizedGlmnet
```

*print a "cv.customizedGlmnet" object*

---

### Description

Print the number of customized training subsets chosen by cross-validation and the number of variables selected in each training subset.

### Usage

```
## S3 method for class 'cv.customizedGlmnet'
print(x, ...)
```

### Arguments

x	a fitted cv.customizedGlmnet object
...	ignored

### Author(s)

Scott Powers, Trevor Hastie, Robert Tibshirani

### See Also

[print, cv.customizedGlmnet](#)

### Examples

```
require(glmnet)

# Simulate synthetic data

n = m = 150
p = 50
q = 5
K = 3
sigmaC = 10
sigmaX = sigmaY = 1
set.seed(5914)
```

```

beta = matrix(0, nrow = p, ncol = K)
for (k in 1:K) beta[sample(1:p, q), k] = 1
c = matrix(rnorm(K*p, 0, sigmaC), K, p)
eta = rnorm(K)
pi = (exp(eta)+1)/sum(exp(eta)+1)
z = t(rmultinom(m + n, 1, pi))
x = crossprod(t(z), c) + matrix(rnorm((m + n)*p, 0, sigmaX), m + n, p)
y = rowSums(z*(crossprod(t(x), beta))) + rnorm(m + n, 0, sigmaY)

x.train = x[1:n, ]
y.train = y[1:n]
x.test = x[n + 1:m, ]
y.test = y[n + 1:m]
foldid = sample(rep(1:10, length = nrow(x.train)))

# Example 1: Use clustering to fit the customized training model to training
# and test data with no predefined test-set blocks

fit1 = cv.customizedGlmnet(x.train, y.train, x.test, Gs = c(1, 2, 3, 5),
  family = "gaussian", foldid = foldid)

# Print the optimal number of groups and value of lambda:
fit1$G.min
fit1$lambda.min

# Print the customized training model fit:
fit1

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit1))^2)

# Plot nonzero coefficients by group:
plot(fit1)

# Example 2: If the test set has predefined blocks, use these blocks to define
# the customized training sets, instead of using clustering.
foldid = apply(z == 1, 1, which)[1:n]
group.id = apply(z == 1, 1, which)[n + 1:m]

fit2 = cv.customizedGlmnet(x.train, y.train, x.test, group.id, foldid = foldid)

# Print the optimal value of lambda:
fit2$lambda.min

# Print the customized training model fit:
fit2

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit2))^2)

```



```
# Plot nonzero coefficients by group:
plot(fit2)

# Example 3: If there is no test set, but the training set is organized into
# blocks, you can do cross validation with these blocks as the basis for the
# customized training sets.

fit3 = cv.customizedGlmnet(x.train, y.train, foldid = foldid)

# Print the optimal value of lambda:
fit3$lambda.min

# Print the customized training model fit:
fit3

# Compute test error using the predict function:
mean((y[n + 1:m] - predict(fit3))^2)

# Plot nonzero coefficients by group:
plot(fit3)
```

---

Vowel

*Vowel Recognition*

---

### **Description**

Speaker independent recognition of the eleven steady state vowels of British English using a specified training set of lpc derived log area ratios.

### **Format**

A data frame with 990 observations on the following 12 variables.

y Class label indicating vowel spoken  
subset a factor with levels test train  
x.1 a numeric vector  
x.2 a numeric vector  
x.3 a numeric vector  
x.4 a numeric vector  
x.5 a numeric vector  
x.6 a numeric vector  
x.7 a numeric vector  
x.8 a numeric vector  
x.9 a numeric vector  
x.10 a numeric vector

### Details

The speech signals were low pass filtered at 4.7kHz and then digitised to 12 bits with a 10kHz sampling rate. Twelfth order linear predictive analysis was carried out on six 512 sample Hamming windowed segments from the steady part of the vowel. The reflection coefficients were used to calculate 10 log area parameters, giving a 10 dimensional input space. For a general introduction to speech processing and an explanation of this technique see Rabiner and Schafer [RabinerSchafer78].

Each speaker thus yielded six frames of speech from eleven vowels. This gave 528 frames from the eight speakers used to train the networks and 462 frames from the seven speakers used to test the networks.

The eleven vowels, along with words demonstrating their sound, are: i (heed) I (hid) E (head) A (had) a: (hard) Y (hud) O (hod) C: (hoard) U (hood) u: (who'd) ɜ: (heard)

### Source

<https://archive.ics.uci.edu/ml/machine-learning-databases/undocumented/connectionist-bench/vowel/>

### References

D. H. Deterding, 1989, University of Cambridge, "Speaker Normalisation for Automatic Speech Recognition", submitted for PhD.

### Examples

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summary(Vowel)
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