

# Package: classify (via r-universe)

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**Title** Explore Classification Models in High Dimensions

**Version** 0.4.3

**Description** Given  $p$ -dimensional training data containing  $d$  groups (the design space), a classification algorithm (classifier) predicts which group new data belongs to. Generally the input to these algorithms is high dimensional, and the boundaries between groups will be high dimensional and perhaps non-linear. This package implements methods for understanding the division of space between the groups.

**License** MIT + file LICENSE

**URL** <http://had.co.nz/classify>

**Imports** class, plyr, stats

**Suggests** e1071, MASS, rpart

**Encoding** UTF-8

**LazyData** true

**RoxygenNote** 7.3.3

**NeedsCompilation** no

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advantage	<i>Calculate the advantage the most likely class has over the next most likely.</i>
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### Description

This is used to identify the boundaries between classification regions. Points with low (close to 0) advantage are likely to be near boundaries.

### Usage

```
advantage(post)
```

### Arguments

post	matrix of posterior probabilities
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classify	<i>Classify provides a convenient method to fit a classification function and then explore the results in the original high dimensional space.</i>
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### Description

This is a convenient function to fit a classification function and then explore the results using GGobi. You can also do this in two separate steps using the classification function and then [explore](#).

### Usage

```
classify(
  data,
  model,
  classifier,
  ...,
  n = 10000,
  method = "nonaligned",
  type = "range"
)
```

**Arguments**

data	Data set use for classification
model	Classification formula, usually of the form response ~ predictors
classifier	Function to use for the classification, eg. <a href="#">lda</a>
...	Other arguments passed to classification function. For example. if you use <a href="#">svm</a> you need to use probability = TRUE so that posterior probabilities can be retrieved.
n	Number of points to simulate. To maintain the illusion of a filled solid this needs to increase with dimension. 10,000 points seems adequate for up to four of five dimensions, but if you have more predictors than that, you will need to increase this number.
method	method to simulate points: grid, random or nonaligned (default). See <a href="#">simvar</a> for more details on the methods used.
type	type of scaling to apply to data. Defaults to common range. See <a href="#">rescaler</a> for more details.

**Details**

By default in GGobi, points that are not on the boundary (ie. that have an advantage greater than the 5% percentile) are hidden. To show them, switch to brush mode and choose include shadowed points from the brush menu on the plot window. You can then brush them yourself to explore how the certainty of classification varies throughout the space

Special notes:

- You should make sure the response variable is a factor
- For SVM, make sure to include probability = TRUE in the arguments to classify

**See Also**

[explore](#), <http://had.co.nz/classify>

**Examples**

```
data(kyphosis, package = "rpart")
library(MASS)
classify(kyphosis, Kyphosis ~ . , lda)
classify(kyphosis, Kyphosis ~ . , qda)
classify(kyphosis, Kyphosis ~ . , glm, family="binomial")
classify(kyphosis, Kyphosis ~ . , knnf, k=3)

library(rpart)
classify(kyphosis, Kyphosis ~ . , rpart)

if (require("e1071")) {
  classify(kyphosis, Kyphosis ~ . , svm, probability=TRUE)
  classify(kyphosis, Kyphosis ~ . , svm, probability=TRUE, kernel="linear")
}
```

```

classify(kyphosis, Kyphosis ~ . , best.svm, probability=TRUE,
         kernel="linear")

# Also can use explore directly
bsvm <- best.svm(Species~., data = iris, gamma = 2^(-1:1),
               cost = 2^(2:+ 4), probability=TRUE)
explore(bsvm, iris)
}

```

---

explore

*Default method for exploring objects*

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### Description

The default method currently works for classification functions.

### Usage

```
explore(model, data, n = 10000, method = "nonaligned", advantage = TRUE, ...)
```

### Arguments

model	classification object
data	data set used with classifier
n	number of points to generate when searching for boundaries
method	method to generate points, see <a href="#">generate_data</a>
advantage	only display boundaries
...	other arguments not currently used

### Details

It generates a data set filling the design space, finds class boundaries (if desired) and then displays in a new ggobi instance.

### Value

A [invisible](#) data frame of class `classify` that contains all the simulated and true data. This can be saved and then printed later to open with `rggobi`.

### See Also

[generate\\_classification\\_data](#), <http://had.co.nz/classify>

## Examples

```
if (require("e1071")) {  
  bsvm <- best.svm(Species~., data = iris, gamma = 2^(-1:1),  
    cost = 2^(2:+ 4), probability=TRUE)  
  explore(bsvm, iris)  
}
```

---

generate\_classification\_data

*Generate classification data.*

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## Description

Given a model, this function generates points within the range of the data, classifies them, and attempts to locate boundaries by looking at advantage.

## Usage

```
generate_classification_data(model, data, n, method, advantage)
```

## Arguments

model	classification model
data	data set used in model
n	number of points to generate
method	method to use, currently either grid (an evenly spaced grid), random (uniform random distribution across cube), or nonaligned (grid + some random perturbation)
advantage	if TRUE, compute advantage, otherwise don't

## Details

If posterior probabilities of classification are available, then the `advantage()` will be calculated directly. If not, `class::knn()` is used calculate the advantage based on the number of neighbouring points that share the same classification. Because knn is  $O(n^2)$  this method is rather slow for large (>20,000 say) data sets.

By default, the boundary points are identified as those below the 5th-percentile for advantage.

## Value

data.frame of classified data

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generate_data	<i>Generate new data from a data frame.</i>
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**Description**

This method generates new data that fills the range of the supplied datasets.

**Usage**

```
generate_data(data, n = 10000, method = "grid")
```

**Arguments**

data	data frame
n	desired number of new observations
method	method to use, see <a href="#">simvar</a>

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knnf	<i>A wrapper function for <a href="#">knn</a> to allow use with <a href="#">classify</a>.</i>
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**Description**

A wrapper function for [knn](#) to allow use with [classify](#).

**Usage**

```
knnf(formula, data, k = 2)
```

**Arguments**

formula	classification formula
data	training data set
k	number of neighbours to use

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`olives`*Olives*

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**Description**

The olive oil data consists of the percentage composition of 8 fatty acids (palmitic, palmitoleic, stearic, oleic, linoleic, linolenic, arachidic, eicosenoic) found in the lipid fraction of 572 Italian olive oils. There are 9 collection areas, 4 from southern Italy (North and South Apulia, Calabria, Sicily), two from Sardinia (Inland and Coastal) and 3 from northern Italy (Umbria, East and West Liguria).

**Format**

A data frame with 244 rows and 7 variables

**References**

Forina, M. and Armanino, C. and Lanteri, S. and Tiscornia, E., Classification of olive oils from their fatty acid composition, 1983, in Food Research and Data Analysis, edited by Martens, H. and Russwurm Jr, H, pages 189-214.

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`posterior`*Extract posterior group probabilities*

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**Description**

Every classification method seems to provide a slightly different way of retrieving the posterior probability of group membership. This function provides a common interface to all of them

**Usage**

```
posterior(model, data)
```

**Arguments**

<code>model</code>	model object
<code>data</code>	data set used in model

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simvar	<i>Simulate observations from a vector</i>
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**Description**

Given a vector of data this function will simulate data that could have come from that vector.

**Usage**

```
simvar(x, n = 10, method = "grid")
```

**Arguments**

x	data vector
n	desired number of points (will not always be achieved)
method	grid simulation method. See details.

**Details**

There are three methods to choose from:

- nonaligned (default): grid + some random perturbation
- grid: grid of evenly spaced observations. If a factor, all levels in a factor will be used, regardless of n
- random: a random uniform sample from the range of the variable

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variables	<i>Extract predictor and response variables for a model object.</i>
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**Description**

Due to the way that most model objects are stored, you also need to supply the data set you used with the original data set. It currently doesn't support models fitted without using a data argument.

**Usage**

```
variables(model)
```

**Arguments**

model	model object
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**Value**

list containing response and predictor variables

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