Package ‘discrim’

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NeedsCompilation no

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discrim_flexible General Interface for Flexible Discriminant Models

Description

discrim_flexible() is a way to generate a specification of a flexible discriminant model using features created using multivariate adaptive regression splines (MARS).

Usage

discrim_flexible(
  mode = "classification",
  num_terms = NULL,
  prod_degree = NULL,
  prune_method = NULL
)

## S3 method for class 'discrim_flexible'
update(
  object,
  num_terms = NULL,
  prod_degree = NULL,
  prune_method = NULL,
  fresh = FALSE,
  ...
)

Arguments

- `mode` A single character string for the type of model. The only possible value for this model is "classification".
- `num_terms` The number of features that will be retained in the final model, including the intercept.
- `prod_degree` The highest possible interaction degree.
- `prune_method` The pruning method.
- `object` A flexible discriminant model specification.
Discriminative Analysis (FDA) uses the work of Hastie et al. (1994) to create a discriminant model using different feature expansions. For this function, MARS (Friedman, 1991) hinge functions are used to non-linearly model the class boundaries (see example below). The mda and earth packages are needed to fit this model.

The main arguments for the model are:

- `num_terms`: The number of features that will be retained in the final model.
- `prod_degree`: The highest possible degree of interaction between features. A value of 1 indicates an additive model while a value of 2 allows, but does not guarantee, two-way interactions between features.
- `prune_method`: The type of pruning. Possible values are listed in ?earth.

These arguments are converted to their specific names at the time that the model is fit. Other options and argument can be set using `set_engine()`. If left to their defaults here (NULL), the values are taken from the underlying model functions. If parameters need to be modified, `update()` can be used in lieu of recreating the object from scratch.

The model can be created using the `fit()` function using the following `engines`:

- **R**: "earth" (the default)

### Engine Details

Engines may have pre-set default arguments when executing the model fit call. For this type of model, the template of the fit calls are:

```r
discrim_flexible() %>%
  set_engine("earth") %>%
  translate()
```

```r
## Flexible Discriminant Model Specification (classification)
## Computational engine: earth
## Model fit template:
## mda::fda(formula = missing_arg(), data = missing_arg(), method = earth::earth)
```

The standardized parameter names in parsnip can be mapped to their original names in each engine that has main parameters. Each engine typically has a different default value (shown in parentheses) for each parameter.

<table>
<thead>
<tr>
<th>parsnip</th>
<th>earth</th>
</tr>
</thead>
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<tr>
<td><code>num_terms</code></td>
<td><code>n.prune</code> (all created by forward pass)</td>
</tr>
<tr>
<td><code>prod_degree</code></td>
<td><code>degree</code> (1)</td>
</tr>
<tr>
<td><code>prune_method</code></td>
<td><code>p.method</code> (backward)</td>
</tr>
</tbody>
</table>
References


Examples

```r
parabolic_grid <-
  expand.grid(X1 = seq(-5, 5, length = 100),
              X2 = seq(-5, 5, length = 100))

fda_mod <-
  discrim_flexible(num_terms = 3) %>%
  set_engine("earth") %>%
  fit(class ~ ., data = parabolic)

parabolic_grid$fda <-
  predict(fda_mod, parabolic_grid, type = "prob")$.pred_Class1

library(ggplot2)

ggplot(parabolic, aes(x = X1, y = X2)) +
  geom_point(aes(col = class), alpha = .5) +
  geom_contour(data = parabolic_grid, aes(z = fda), col = "black", breaks = .5) +
  theme_bw() +
  theme(legend.position = "top") +
  coord_equal()

model <- discrim_flexible(num_terms = 10)
model
update(model, num_terms = 6)
```

---

discrim_linear

*General Interface for Linear Discriminant Models*

**Description**

discrim_linear() is a way to generate a specification of a linear discriminant analysis (LDA) model before fitting and allows the model to be created using different packages in R.

**Usage**

discrim_linear(mode = "classification", penalty = NULL)

## S3 method for class 'discrim_linear'
update(object, penalty = NULL, fresh = FALSE, ...)

Arguments

- **mode**: A single character string for the type of model. The only possible value for this model is "classification".
- **penalty**: An non-negative number representing the amount of regularization used by some of the engines.
- **object**: A linear discriminant model specification.
- **fresh**: A logical for whether the arguments should be modified in-place of or replaced wholesale.
- **...**: Not used for update().

Details

For `discrim_linear()`, the mode will always be "classification".

The model can be created using the `fit()` function using the following **engines**:

- **R**: "MASS" (the default) or "mda"

The main argument for the model is:

- **penalty**: The total amount of regularization in the model. Note that this only used for the "mda" engine where it is a pure L2 penalty (a.k.a ridge regression).

This argument is converted to its specific names at the time that the model is fit. Other options and argument can be set using `set_engine()`. If left to their defaults here (NULL), the values are taken from the underlying model functions. If parameters need to be modified, `update()` can be used in lieu of recreating the object from scratch.

Engine Details

Engines may have pre-set default arguments when executing the model fit call. For this type of model, the template of the fit calls are:

```r
discrim_linear() %>%
  set_engine("MASS") %>%
  translate()
```

## Linear Discriminant Model Specification (classification)

## Computational engine: MASS

## Model fit template:

```
MASS::lda(formula = missing_arg(), data = missing_arg())
```

```r
discrim_linear() %>%
  set_engine("mda") %>%
  translate()
```
## Linear Discriminant Model Specification (classification)
## Computational engine: mda
## Model fit template:
## mda::fda(formula = missing_arg(), data = missing_arg(), method = mda::gen.ridge, 
## keep.fitted = FALSE)

The standardized parameter names in parsnip can be mapped to their original names in each engine that has main parameters. Each engine typically has a different default value (shown in parentheses) for each parameter.

<table>
<thead>
<tr>
<th>parsnip</th>
<th>mda</th>
</tr>
</thead>
<tbody>
<tr>
<td>penalty</td>
<td>lambda (1)</td>
</tr>
</tbody>
</table>

### Examples

```r
parabolic_grid <-
  expand.grid(X1 = seq(-5, 5, length = 100),
              X2 = seq(-5, 5, length = 100))

lda_mod <-
  discrim_linear(penalty = .1) %>%
  set_engine("mda") %>%
  fit(class ~ ., data = parabolic)

parabolic_grid$lda <-
  predict(lda_mod, parabolic_grid, type = "prob")$.pred_Class1

library(ggplot2)
ggplot(parabolic, aes(x = X1, y = X2)) +
  geom_point(aes(col = class), alpha = .5) +
  geom_contour(data = parabolic_grid, aes(z = lda), col = "black", breaks = .5) +
  theme_bw() +
  theme(legend.position = "top") +
  coord_equal()

model <- discrim_linear(penalty = 0.1)
model
update(model, penalty = 1)
```

---

discrim_regularized  General Interface for Regularized Discriminant Models

### Description

discrim_regularized() is a way to generate a specification of a regularized discriminant analysis (RDA) model before fitting.
Usage

discrim_regularized(
    mode = "classification",
    frac_common_cov = NULL,
    frac_identity = NULL
)

## S3 method for class 'discrim_regularized'
update(
    object,
    frac_common_cov = NULL,
    frac_identity = NULL,
    fresh = FALSE,
    ...
)

Arguments

mode A single character string for the type of model. The only possible value for this model is 'classification'.
frac_common_cov, frac_identity Numeric values between zero and one.
object A linear discriminant model specification.
fresh A logical for whether the arguments should be modified in-place of or replaced wholesale.
... Not used for update().

Details

The model is from Friedman (1989) and can create LDA models, QDA models, and regularized mixtures of the two. It does not conduct feature selection. The main arguments for the model are:

- `frac_common_cov`: The fraction of the regularized covariance matrix that is based on the LDA model (i.e., computed from all classes). A value of 1 is the linear discriminant analysis assumption while a value near zero assumes that there should be separate covariance matrices for each class.
- `frac_identity`: The fraction of the final, class-specific covariance matrix that is the identity matrix.

See klaR::rda() for the equations that define these parameters.

These arguments are converted to their specific names at the time that the model is fit. Other options and argument can be set using set_engine(). If left to their defaults here (NULL), the values are taken from the underlying model functions. If parameters need to be modified, update() can be used in lieu of recreating the object from scratch.

For discrim_regularized(), the mode will always be "classification".
Engine Details

Engines may have pre-set default arguments when executing the model fit call. For this type of model, the template of the fit calls are:

discrim_regularized() %>%
  set_engine("klaR") %>%
  translate()

## Regularized Discriminant Model Specification (classification)
## Computational engine: klaR
## Model fit template:
## klaR::rda(formula = missing_arg(), data = missing_arg())

The standardized parameter names in parsnip can be mapped to their original names in each engine that has main parameters. Each engine typically has a different default value (shown in parentheses) for each parameter.

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<thead>
<tr>
<th>parsnip</th>
<th>klaR</th>
</tr>
</thead>
<tbody>
<tr>
<td>frac_common_cov</td>
<td>lambda (varies)</td>
</tr>
<tr>
<td>frac_identity</td>
<td>gamma (varies)</td>
</tr>
</tbody>
</table>

References


Examples

```r
parabolic_grid <-
  expand.grid(X1 = seq(-5, 5, length = 100),
              X2 = seq(-5, 5, length = 100))

rda_mod <-
  discrim_regularized(frac_common_cov = .5, frac_identity = .5) %>%
  set_engine("klaR") %>%
  fit(class ~ ., data = parabolic)

parabolic_grid$rda <-
  predict(rda_mod, parabolic_grid, type = "prob")$.pred_Class1

library(ggplot2)
ggplot(parabolic, aes(x = X1, y = X2)) +
  geom_point(aes(col = class), alpha = .5) +
  geom_contour(data = parabolic_grid, aes(z = rda), col = "black", breaks = .5) +
  theme_bw() +
  theme(legend.position = "top") +
  coord_equal()
```
model <- discrim_regularized(frac_common_cov = 10)
model
update(model, frac_common_cov = 1)

frac_common_cov  Parameter objects for Regularized Discriminant Models

Description
discrim_regularized() describes the effect of frac_common_cov() and frac_identity(). smoothness() is an alias for the adjust parameter in stats::density().

Usage
frac_common_cov(range = c(0, 1), trans = NULL)
frac_identity(range = c(0, 1), trans = NULL)
smoothness(range = c(0.5, 1.5), trans = NULL)

Arguments
range A two-element vector holding the defaults for the smallest and largest possible values, respectively.
trans A trans object from the scales package, such as scales::log10_trans() or scales::reciprocal_trans(). If not provided, the default is used which matches the units used in range. If no transformation, NULL.

Details
These parameters can modulate a RDA model to go between linear and quadratic class boundaries.

Value
A function with classes "quant_param" and "param"

Examples
frac_common_cov()
**Description**

`naive_Bayes()` is a way to generate a *specification* of a model before fitting and allows the model to be created using different packages in R.

**Usage**

```r
naive_Bayes(mode = "classification", smoothness = NULL, Laplace = NULL)
```

```r
## S3 method for class 'naive_Bayes'
update(object, smoothness = NULL, Laplace = NULL, fresh = FALSE, ...)
```

**Arguments**

- **mode**: A single character string for the type of model. The only possible value for this model is "classification".
- **smoothness**: An non-negative number representing the the relative smoothness of the class boundary. Smaller examples result in model flexible boundaries and larger values generate class boundaries that are less adaptable.
- **Laplace**: A non-negative value for the Laplace correction to smoothing low-frequency counts.
- **object**: A linear discriminant model specification.
- **fresh**: A logical for whether the arguments should be modified in-place of or replaced wholesale.
- **...**: Not used for `update()`.

**Details**

The main arguments for the model are:

- **smoothness**: The total amount of regularization in the model. Note that this only used for the "klaR" engine where it is a pure L2 smoothness (a.k.a ridge regression).
- **Laplace**: Laplace correction for smoothing low-frequency counts.

These arguments are converted to their specific names at the time that the model is fit. Other options and argument can be set using `set_engine()`. If left to their defaults here (NULL), the values are taken from the underlying model functions. If parameters need to be modified, `update()` can be used in lieu of recreating the object from scratch.

For `naive_Bayes()`, the mode will always be "classification".

The model can be created using the `fit()` function using the following **engines**:

- **R**: "klaR" (the default) or "naivebayes"
Engine Details

Engines may have pre-set default arguments when executing the model fit call. For this type of model, the template of the fit calls are:

```r
naive_Bayes() %>%
  set_engine("klaR") %>%
  translate()

## Naive Bayes Model Specification (classification)
## Computational engine: klaR
## Model fit template:
## discrim::klar_bayes_wrapper(x = missing_arg(), y = missing_arg(),
## usekernel = TRUE)
```

```r
naive_Bayes() %>%
  set_engine("naivebayes") %>%
  translate()

## Naive Bayes Model Specification (classification)
## Computational engine: naivebayes
## Model fit template:
## naivebayes::naive_bayes(x = missing_arg(), y = missing_arg(),
## usekernel = TRUE)
```

The standardized parameter names in parsnip can be mapped to their original names in each engine that has main parameters. Each engine typically has a different default value (shown in parentheses) for each parameter.

<table>
<thead>
<tr>
<th>parsnip</th>
<th>klaR</th>
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</tr>
</thead>
<tbody>
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<td>smoothness</td>
<td>adjust (1)</td>
<td>adjust (1)</td>
</tr>
<tr>
<td>Laplace</td>
<td>fL (0)</td>
<td>laplace (0)</td>
</tr>
</tbody>
</table>

Note that `usekernel` is always set to `TRUE` for the `klaR` engine.

Examples

```r
parabolic_grid <-
  expand.grid(X1 = seq(-5, 5, length = 100),
              X2 = seq(-5, 5, length = 100))

nb_mod <-
  naive_Bayes(smoothness = .8) %>%
  set_engine("klaR") %>%
```
```r
library(ggplot2)
ggplot(parabolic, aes(x = X1, y = X2)) +
  geom_point(aes(col = class), alpha = .5) +
  geom_contour(data = parabolic_grid, aes(z = nb), col = "black", breaks = .5) +
  theme_bw() +
  theme(legend.position = "top") +
  coord_equal()
```

```r
model <- naive_Bayes(smoothness = 0.1)
model
update(model, smoothness = 1)
update(model, smoothness = 1, fresh = TRUE)
```

---

**parabolic**

*Parabolic class boundary data*

**Description**

Parabolic class boundary data

**Details**

These data were simulated. There are two correlated predictors and two classes in the factor outcome.

**Value**

parabolic a data frame

**Examples**

data(parabolic)

library(ggplot2)
ggplot(parabolic, aes(x = X1, y = X2, col = class)) +
  geom_point(alpha = .5) +
  theme_bw()
```
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