

Package ‘baguette’

April 5, 2023

Title Efficient Model Functions for Bagging

Version 1.0.1

Description Tree- and rule-based models can be bagged (<[doi:10.1007/BF00058655](https://doi.org/10.1007/BF00058655)>) using this package and their predictions equations are stored in an efficient format to reduce the model objects size and speed.

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URL <https://baguette.tidymodels.org>,
<https://github.com/tidymodels/baguette>

BugReports <https://github.com/tidymodels/baguette/issues>

Depends parsnip (>= 1.0.0), R (>= 3.4)

Imports butcher, C50, dials, dplyr, earth, furrr, generics, hardhat (>= 1.1.0), magrittr, purrr, rlang, rpart, rsample, tibble, tidyr, utils, withr

Suggests AmesHousing, covr, modeldata, nnet, recipes, rmarkdown, spelling, testthat (>= 3.0.0), yardstick

Config/Needs/website tidyverse/tidytemplate

Config/testthat/edition 3

Encoding UTF-8

Language en-US

RoxygenNote 7.2.3

NeedsCompilation no

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Repository CRAN

Date/Publication 2023-04-04 23:00:02 UTC

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|--------|--------------------------|
| bagger | <i>Bagging functions</i> |
|--------|--------------------------|

Description

General suite of bagging functions for several models.

Usage

```
bagger(x, ...)
```

```
## Default S3 method:
bagger(x, ...)
```

```
## S3 method for class 'data.frame'
bagger(
  x,
  y,
  weights = NULL,
  base_model = "CART",
  times = 11L,
  control = control_bag(),
  cost = NULL,
  ...
)
```

```
## S3 method for class 'matrix'
bagger(
  x,
  y,
  weights = NULL,
  base_model = "CART",
  times = 11L,
  control = control_bag(),
  cost = NULL,
  ...
)
```

```

## S3 method for class 'formula'
bagger(
  formula,
  data,
  weights = NULL,
  base_model = "CART",
  times = 11L,
  control = control_bag(),
  cost = NULL,
  ...
)

## S3 method for class 'recipe'
bagger(
  x,
  data,
  base_model = "CART",
  times = 11L,
  control = control_bag(),
  cost = NULL,
  ...
)

```

Arguments

| | |
|-------------------------|--|
| <code>x</code> | A data frame, matrix, or recipe (depending on the method being used). |
| <code>...</code> | Optional arguments to pass to the base model function. |
| <code>y</code> | A numeric or factor vector of outcomes. Categorical outcomes (i.e classes) should be represented as factors, not integers. |
| <code>weights</code> | A numeric vector of non-negative case weights. These values are not used during bootstrap resampling. |
| <code>base_model</code> | A single character value for the model being bagged. Possible values are "CART", "MARS", "nnet", and "C5.0" (classification only). |
| <code>times</code> | A single integer greater than 1 for the maximum number of bootstrap samples/ensemble members (some model fits might fail). |
| <code>control</code> | A list of options generated by <code>control_bag()</code> . |
| <code>cost</code> | A non-negative scale (for two class problems) or a cost matrix. |
| <code>formula</code> | An object of class "formula" (or one that can be coerced to that class): a symbolic description of the model to be fitted. Note that this package does not support multivariate outcomes and that, if some predictors are factors, dummy variables will <i>not</i> be created unless by the underlying model function. |
| <code>data</code> | A data frame containing the variables used in the formula or recipe. |

Details

`bagger()` fits separate models to bootstrap samples. The prediction function for each model object

is encoded in an R expression and the original model object is discarded. When making predictions, each prediction formula is evaluated on the new data and aggregated using the mean.

Variable importance scores are calculated using implementations in each package. When requested, the results are in a tibble with column names `term` (the predictor), `value` (the importance score), and `used` (the percentage of times that the variable was in the prediction equation).

The models can be fit in parallel using the **future** package. To enable parallelism, use the `future::plan()` function to declare *how* the computations should be distributed. Note that this will almost certainly multiply the memory requirements required to fit the models.

For neural networks, variable importance is calculated using the method of Garson described in Gevrey *et al* (2003)

References

Gevrey, M., Dimopoulos, I., and Lek, S. (2003). Review and comparison of methods to study the contribution of variables in artificial neural network models. *Ecological Modelling*, 160(3), 249-264.

Examples

```
library(recipes)
library(dplyr)

data(biomass, package = "modeldata")

biomass_tr <-
  biomass %>%
  dplyr::filter(dataset == "Training") %>%
  dplyr::select(-dataset, -sample)

biomass_te <-
  biomass %>%
  dplyr::filter(dataset == "Testing") %>%
  dplyr::select(-dataset, -sample)

# -----

ctrl <- control_bag(var_imp = TRUE)

# -----

# `times` is low to make the examples run faster

set.seed(7687)
mars_bag <- bagger(x = biomass_tr[, -6], y = biomass_tr$HHV,
                  base_model = "MARS", times = 5, control = ctrl)
mars_bag
var_imp(mars_bag)

set.seed(7687)
cart_bag <- bagger(x = biomass_tr[, -6], y = biomass_tr$HHV,
                  base_model = "CART", times = 5, control = ctrl)
```

```

cart_bag

# -----
# Other interfaces

# Recipes can be used
biomass_rec <-
  recipe(HHV ~ ., data = biomass_tr) %>%
  step_pca(all_predictors())

set.seed(7687)
cart_pca_bag <- bagger(biomass_rec, data = biomass_tr, base_model = "CART",
                      times = 5, control = ctrl)

cart_pca_bag

# Using formulas
mars_bag <- bagger(HHV ~ ., data = biomass_tr, base_model = "MARS", times = 5,
                  control = ctrl)

mars_bag

```

class_cost

Cost parameter for minority class

Description

Used in `bag_treer()`.

Usage

```
class_cost(range = c(0, 5), trans = NULL)
```

Arguments

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

Details

This parameter reflects the cost of a misclassified sample relative to a baseline cost of 1.0. For example, if the first level of an outcome factor occurred rarely, it might help if this parameter were set to values greater than 1.0. If the second level of the outcome factor is in the minority, values less than 1.0 would cause the model to emphasize the minority class more than the majority class.

Examples

```
class_cost()
```

control_bag

*Controlling the bagging process***Description**

control_bag() can set options for ancillary aspects of the bagging process.

Usage

```
control_bag(
  var_imp = TRUE,
  allow_parallel = TRUE,
  sampling = "none",
  reduce = TRUE,
  extract = NULL
)
```

Arguments

| | |
|----------------|--|
| var_imp | A single logical: should variable importance scores be calculated? |
| allow_parallel | A single logical: should the model fits be done in parallel (even if a parallel plan() has been created)? |
| sampling | Either "none" or "down". For classification only. The training data, after bootstrapping, will be sampled down within each class (with replacement) to the size of the smallest class. |
| reduce | Should models be modified to reduce their size on disk? |
| extract | A function (or NULL) that can extract model-related aspects of each ensemble member. See Details and example below. |

Details

Any arbitrary item can be saved from the model object (including the model object itself) using the extract argument, which should be a function with arguments x (for the model object), and ... The results of this function are saved into a list column called extras (see the example below).

Value

A list.

Examples

```
# Extracting model components

num_term_nodes <- function(x, ...) {
  tibble::tibble(num_nodes = sum(x$frame$var == "<leaf>"))
}
```

```

set.seed(7687)
with_extras <- bagger(mpg ~ ., data = mtcars,
                      base_model = "CART", times = 5,
                      control = control_bag(extract = num_term_nodes))

dplyr::bind_rows(with_extras$model_df$extras)

```

predict.bagger *Predictions from a bagged model*

Description

The `predict()` function computes predictions from each of the models in the ensembles and returns a single aggregated value for each sample in `new_data`.

Usage

```

## S3 method for class 'bagger'
predict(object, new_data, type = NULL, ...)

```

Arguments

| | |
|-----------------------|---|
| <code>object</code> | An object generated by <code>bagger()</code> . |
| <code>new_data</code> | A data frame of predictors. If a recipe or formula were originally used, the original data should be passed here instead of a preprocessed version. |
| <code>type</code> | A single character value for the type of predictions. For regression models, <code>type = 'numeric'</code> is valid and <code>'class'</code> and <code>'prob'</code> are valid for classification models. |
| <code>...</code> | Not currently used. |

Examples

```

data(airquality)

set.seed(7687)
mars_bag <- bagger(Ozone ~ ., data = airquality, base_model = "MARS", times = 5)
predict(mars_bag, new_data = airquality[, -1])

```

| | |
|----------------|--|
| var_imp.bagger | <i>Obtain variable importance scores</i> |
|----------------|--|

Description

Obtain variable importance scores

Usage

```
## S3 method for class 'bagger'
var_imp(object, ...)
```

Arguments

| | |
|--------|---------------------|
| object | An object. |
| ... | Not currently used. |

Details

baguette can compute different variable importance scores for each model in the ensemble. The `var_imp()` function returns the average importance score for each model. Additionally, the function returns the number of times that each predictor is included in the final prediction equation.

Specific methods used by the models are:

CART: The model accumulates the improvement of the model that occurs when a predictor is used in a split. These values are taken from the `rpart` object. See `rpart::rpart.object()`.

MARS: MARS models include a backwards elimination feature selection routine that looks at reductions in the generalized cross-validation (GCV) estimate of error. The `earth()` function tracks the changes in model statistics, such as the GCV, for each predictor and accumulates the reduction in the statistic when each predictor's feature is added to the model. This total reduction is used as the variable importance measure. If a predictor was never used in any of the MARS basis functions in the final model (after pruning), it has an importance value of zero. `baguette` wraps `earth::evimp()`.

C5.0: `C5.0` measures predictor importance by determining the percentage of training set samples that fall into all the terminal nodes after the split. For example, the predictor in the first split automatically has an importance measurement of 100 percent since all samples are affected by this split. Other predictors may be used frequently in splits, but if the terminal nodes cover only a handful of training set samples, the importance scores may be close to zero.

Note that the value column that is the average of the importance scores from each model. The divisor of this average (and the corresponding standard error) is the number of models (as opposed to the number of models that used the predictor). This means that the importance scores for a predictor that was not used in the model has an implicit zero importance.

Value

A tibble with columns for `term` (the predictor), `value` (the mean importance score), `std.error` (the standard error), and `used` (the occurrences of the predictors).

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