

Package ‘embed’

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Title Extra Recipes for Encoding Predictors

Version 1.2.2

Description Predictors can be converted to one or more numeric representations using a variety of methods. Effect encodings using simple generalized linear models <[doi:10.48550/arXiv.1611.09477](https://doi.org/10.48550/arXiv.1611.09477)> or nonlinear models <[doi:10.48550/arXiv.1604.06737](https://doi.org/10.48550/arXiv.1604.06737)> can be used. There are also functions for dimension reduction and other approaches.

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

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

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add_woe	<i>Add WoE in a data frame</i>
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Description

A tidyverse friendly way to plug WoE versions of a set of predictor variables against a given binary outcome.

Usage

```
add_woe(.data, outcome, ..., dictionary = NULL, prefix = "woe")
```

Arguments

.data	A <code>tbl</code> . The <code>data.frame</code> to plug the new woe version columns.
outcome	The bare name of the outcome variable.
...	Bare names of predictor variables, passed as you would pass variables to <code>dplyr::select()</code> . This means that you can use all the helpers like <code>starts_with()</code> and <code>matches()</code> .
dictionary	A <code>tbl</code> . If <code>NULL</code> the function will build a dictionary with those variables passed to <code>....</code> . You can pass a custom dictionary too, see dictionary() for details.
prefix	A character string that will be the prefix to the resulting new variables.

Details

You can pass a custom dictionary to `add_woe()`. It must have the exactly the same structure of the output of `dictionary()`. One easy way to do this is to tweak a output returned from it.

Value

A tibble with the original columns of `.data` plus the woe columns wanted.

Examples

```
mtcars |> add_woe("am", cyl, gear:carb)
```

dictionary	<i>Weight of evidence dictionary</i>
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Description

Builds the woe dictionary of a set of predictor variables upon a given binary outcome. Convenient to make a woe version of the given set of predictor variables and also to allow one to tweak some woe values by hand.

Usage

```
dictionary(.data, outcome, ..., Laplace = 1e-06)
```

Arguments

<code>.data</code>	A <code>tbl</code> . The <code>data.frame</code> where the variables come from.
<code>outcome</code>	The bare name of the outcome variable with exactly 2 distinct values.
<code>...</code>	bare names of predictor variables or selectors accepted by <code>dplyr::select()</code> .
<code>Laplace</code>	Default to <code>1e-6</code> . The pseudocount parameter of the Laplace Smoothing estimator. Value to avoid <code>-Inf/Inf</code> from predictor category with only one outcome class. Set to 0 to allow <code>Inf/-Inf</code> .

Details

You can pass a custom dictionary to `step_woe()`. It must have the exactly the same structure of the output of `dictionary()`. One easy way to do this is by tweaking an output returned from it.

Value

a tibble with summaries and woe for every given predictor variable stacked up.

References

Kullback, S. (1959). *Information Theory and Statistics*. Wiley, New York.

Hastie, T., Tibshirani, R. and Friedman, J. (1986). *Elements of Statistical Learning*, Second Edition, Springer, 2009.

Good, I. J. (1985), "Weight of evidence: A brief survey", *Bayesian Statistics*, 2, pp.249-270.

Examples

```
mtcars |> dictionary("am", cyl, gear:carb)
```

solubility	<i>Compound solubility data</i>
------------	---------------------------------

Description

Compound solubility data

Details

Tetko et al. (2001) and Huuskonen (2000) investigated a set of compounds with corresponding experimental solubility values using complex sets of descriptors. They used linear regression and neural network models to estimate the relationship between chemical structure and solubility. For our analyses, we will use 1267 compounds and a set of more understandable descriptors that fall into one of three groups: 208 binary "fingerprints" that indicate the presence or absence of a particular chemical sub-structure, 16 count descriptors (such as the number of bonds or the number of Bromine atoms) and 4 continuous descriptors (such as molecular weight or surface area).

Value

solubility a data frame

Source

Tetko, I., Tanchuk, V., Kasheva, T., and Villa, A. (2001). Estimation of aqueous solubility of chemical compounds using E-state indices. *Journal of Chemical Information and Computer Sciences*, 41(6), 1488-1493.

Huuskonen, J. (2000). Estimation of aqueous solubility for a diverse set of organic compounds based on molecular topology. *Journal of Chemical Information and Computer Sciences*, 40(3), 773-777.

Examples

```
data(solubility)
str(solubility)
```

step_collapse_cart *Supervised Collapsing of Factor Levels*

Description

`step_collapse_cart()` creates a *specification* of a recipe step that can collapse factor levels into a smaller set using a supervised tree.

Usage

```
step_collapse_cart(  
  recipe,  
  ...,  
  role = NA,  
  trained = FALSE,  
  outcome = NULL,  
  cost_complexity = 1e-04,  
  min_n = 5,  
  results = NULL,  
  skip = FALSE,  
  id = rand_id("step_collapse_cart")  
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose which variables are affected by the step. See recipes::selections for more details. For the <code>tidy</code> method, these are not currently used.
<code>role</code>	Not used by this step since no new variables are created.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>outcome</code>	A call to <code>vars</code> to specify which variable is used as the outcome to train CART models in order to pool factor levels.
<code>cost_complexity</code>	A non-negative value that regulates the complexity of the tree when pruning occurs. Values near 0.1 usually correspond to a tree with a single splits. Values of zero correspond to unpruned tree.
<code>min_n</code>	An integer for how many data points are required to make further splits during the tree growing process. Larger values correspond to less complex trees.
<code>results</code>	A list of results to convert to new factor levels.
<code>skip</code>	A logical. Should the step be skipped when the recipe is baked by recipes::bake ? While all operations are baked when recipes::prep is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations

id	A character string that is unique to this step to identify it.
-----------	--

Details

This step uses a CART tree (classification or regression) to group the existing factor levels into a potentially smaller set. It changes the levels in the factor predictor (and the `tidy()` method can be used to understand the translation).

There are a few different ways that the step will not be able to collapse levels. If the model fails or, if the results have each level being in its own split, the original factor levels are retained. There are also cases where there is "no admissible split" which means that the model could not find any signal in the data.

Value

An updated recipe step.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `old`, `new`, and `id`:

terms character, the selectors or variables selected
old character, the old levels
new character, the new levels
id character, id of this step

Case weights

The underlying operation does not allow for case weights.

See Also

[step_collapse_stringdist\(\)](#)

Examples

```
data(ames, package = "modeldata")
ames$Sale_Price <- log10(ames$Sale_Price)

rec <-
  recipe(Sale_Price ~ ., data = ames) |>
  step_collapse_cart(
    Sale_Type, Garage_Type, Neighborhood,
    outcome = vars(Sale_Price)
  ) |>
  prep()
tidy(rec, number = 1)
```

step_collapse_stringdist
collapse factor levels using stringdist

Description

`step_collapse_stringdist()` creates a *specification* of a recipe step that will collapse factor levels that have a low stringdist between them.

Usage

```
step_collapse_stringdist(  
  recipe,  
  ...,  
  role = NA,  
  trained = FALSE,  
  distance = NULL,  
  method = "osa",  
  options = list(),  
  results = NULL,  
  columns = NULL,  
  skip = FALSE,  
  id = rand_id("collapse_stringdist")  
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose which variables are affected by the step. See recipes::selections for more details. For the <code>tidy</code> method, these are not currently used.
<code>role</code>	Not used by this step since no new variables are created.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>distance</code>	Integer, value to determine which strings should be collapsed with which. The value is being used inclusive, so 2 will collapse levels that have a string distance between them of 2 or lower.
<code>method</code>	Character, method for distance calculation. The default is "osa", see stringdist::stringdist-metrics .
<code>options</code>	List, other arguments passed to stringdist::stringdistmatrix() such as <code>weight</code> , <code>q</code> , <code>p</code> , and <code>bt</code> , that are used for different values of <code>method</code> .
<code>results</code>	A list denoting the way the labels should be collapses is stored here once this preprocessing step has be trained by recipes::prep .
<code>columns</code>	A character string of variable names that will be populated (eventually) by the <code>terms</code> argument.

skip	A logical. Should the step be skipped when the recipe is baked by <code>bake()</code> ? While all operations are baked when <code>prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations.
id	A character string that is unique to this step to identify it.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the columns that will be affected) and `base`.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `from`, `to`, and `id`:

terms character, the selectors or variables selected
from character, the old levels
to character, the new levels
id character, id of this step

Case weights

The underlying operation does not allow for case weights.

See Also

[step_collapse_cart\(\)](#)

Examples

```
library(recipes)
library(tibble)
data0 <- tibble(
  x1 = c("a", "b", "d", "e", "sfgsfgsd", "hjhgfjgr"),
  x2 = c("ak", "b", "djj", "e", "hjhgfjgr", "hjhgfjgr")
)

rec <- recipe(~., data = data0) |>
  step_collapse_stringdist(all_predictors(), distance = 1) |>
  prep()

rec |>
  bake(new_data = NULL)

tidy(rec, 1)

rec <- recipe(~., data = data0) |>
  step_collapse_stringdist(all_predictors(), distance = 2) |>
  prep()
```

```
rec |>
  bake(new_data = NULL)

tidy(rec, 1)
```

step_discretize_cart *Discretize numeric variables with CART*

Description

`step_discretize_cart()` creates a *specification* of a recipe step that will discretize numeric data (e.g. integers or doubles) into bins in a supervised way using a CART model.

Usage

```
step_discretize_cart(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  cost_complexity = 0.01,
  tree_depth = 10,
  min_n = 20,
  rules = NULL,
  skip = FALSE,
  id = rand_id("discretize_cart")
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose which variables are affected by the step. See recipes::selections for more details.
<code>role</code>	Defaults to "predictor".
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>outcome</code>	A call to <code>vars</code> to specify which variable is used as the outcome to train CART models in order to discretize explanatory variables.
<code>cost_complexity</code>	The regularization parameter. Any split that does not decrease the overall lack of fit by a factor of <code>cost_complexity</code> is not attempted. Corresponds to <code>cp</code> in rpart::rpart() . Defaults to 0.01.

tree_depth	The <i>maximum</i> depth in the final tree. Corresponds to <code>maxdepth</code> in <code>rpart::rpart()</code> . Defaults to 10.
min_n	The number of data points in a node required to continue splitting. Corresponds to <code>minsplit</code> in <code>rpart::rpart()</code> . Defaults to 20.
rules	The splitting rules of the best CART tree to retain for each variable. If length zero, splitting could not be used on that column.
skip	A logical. Should the step be skipped when the recipe is baked by <code>recipes::bake()</code> ? While all operations are baked when <code>recipes::prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

`step_discretize_cart()` creates non-uniform bins from numerical variables by utilizing the information about the outcome variable and applying a CART model.

The best selection of buckets for each variable is selected using the standard cost-complexity pruning of CART, which makes this discretization method resistant to overfitting.

This step requires the **rpart** package. If not installed, the step will stop with a note about installing the package.

Note that the original data will be replaced with the new bins.

Value

An updated version of `recipe` with the new step added to the sequence of any existing operations.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `value`, and `id`:

terms character, the selectors or variables selected
value numeric, location of the splits
id character, id of this step

Tuning Parameters

This step has 3 tuning parameters:

- `cost_complexity`: Cost-Complexity Parameter (type: double, default: 0.01)
- `tree_depth`: Tree Depth (type: integer, default: 10)
- `min_n`: Minimal Node Size (type: integer, default: 20)

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in `recipes::case_weights` and the examples on tidymodels.org.

See Also

[recipes::step_discretize\(\)](#), [step_discretize_xgb\(\)](#)

Examples

```
library(modeldata)
data(ad_data)
library(rsample)

split <- initial_split(ad_data, strata = "Class")

ad_data_tr <- training(split)
ad_data_te <- testing(split)

cart_rec <-
  recipe(Class ~ ., data = ad_data_tr) |>
  step_discretize_cart(
    tau, age, p_tau, Ab_42,
    outcome = "Class", id = "cart splits"
  )

cart_rec <- prep(cart_rec, training = ad_data_tr)

# The splits:
tidy(cart_rec, id = "cart splits")

bake(cart_rec, ad_data_te, tau)
```

step_discretize_xgb *Discretize numeric variables with XgBoost*

Description

`step_discretize_xgb()` creates a *specification* of a recipe step that will discretize numeric data (e.g. integers or doubles) into bins in a supervised way using an XgBoost model.

Usage

```
step_discretize_xgb(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  sample_val = 0.2,
  learn_rate = 0.3,
  num_breaks = 10,
  tree_depth = 1,
```

```

  min_n = 5,
  rules = NULL,
  skip = FALSE,
  id = rand_id("discretize_xgb")
)

```

Arguments

recipe	A recipe object. The step will be added to the sequence of operations for this recipe.
...	One or more selector functions to choose which variables are affected by the step. See recipes::selections for more details.
role	Defaults to "predictor".
trained	A logical to indicate if the quantities for preprocessing have been estimated.
outcome	A call to <code>vars</code> to specify which variable is used as the outcome to train XgBoost models in order to discretize explanatory variables.
sample_val	Share of data used for validation (with early stopping) of the learned splits (the rest is used for training). Defaults to 0.20.
learn_rate	The rate at which the boosting algorithm adapts from iteration-to-iteration. Corresponds to <code>eta</code> in the xgboost package. Defaults to 0.3.
num_breaks	The <i>maximum</i> number of discrete bins to bucket continuous features. Corresponds to <code>max_bin</code> in the xgboost package. Defaults to 10.
tree_depth	The maximum depth of the tree (i.e. number of splits). Corresponds to <code>max_depth</code> in the xgboost package. Defaults to 1.
min_n	The minimum number of instances needed to be in each node. Corresponds to <code>min_child_weight</code> in the xgboost package. Defaults to 5.
rules	The splitting rules of the best XgBoost tree to retain for each variable.
skip	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

`step_discretize_xgb()` creates non-uniform bins from numerical variables by utilizing the information about the outcome variable and applying the **xgboost** model. It is advised to impute missing values before this step. This step is intended to be used particularly with linear models because thanks to creating non-uniform bins it becomes easier to learn non-linear patterns from the data.

The best selection of buckets for each variable is selected using an internal early stopping scheme implemented in the **xgboost** package, which makes this discretization method prone to overfitting.

The pre-defined values of the underlying **xgboost** learns good and reasonably complex results. However, if one wishes to tune them the recommended path would be to first start with changing the value of `num_breaks` to e.g.: 20 or 30. If that doesn't give satisfactory results one could experiment

with modifying the `tree_depth` or `min_n` parameters. Note that it is not recommended to tune `learn_rate` simultaneously with other parameters.

This step requires the **xgboost** package. If not installed, the step will stop with a note about installing the package.

Note that the original data will be replaced with the new bins.

Value

An updated version of `recipe` with the new step added to the sequence of any existing operations.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `value`, and `id`:

`terms` character, the selectors or variables selected

`value` numeric, location of the splits

`id` character, id of this step

Tuning Parameters

This step has 5 tuning parameters:

- `sample_val`: Proportion of data for validation (type: double, default: 0.2)
- `learn_rate`: Learning Rate (type: double, default: 0.3)
- `num_breaks`: Number of Cut Points (type: integer, default: 10)
- `tree_depth`: Tree Depth (type: integer, default: 1)
- `min_n`: Minimal Node Size (type: integer, default: 5)

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in `recipes::case_weights` and the examples on tidymodels.org.

See Also

`recipes::step_discretize()`, `step_discretize_cart()`

Examples

```
library(rsample)
library(recipes)
data(credit_data, package = "modeldata")

set.seed(1234)
split <- initial_split(credit_data[1:1000, ], strata = "Status")

credit_data_tr <- training(split)
credit_data_te <- testing(split)
```

```

xgb_rec <-
  recipe(Status ~ Income + Assets, data = credit_data_tr) |>
  step_impute_median(Income, Assets) |>
  step_discretize_xgb(Income, Assets, outcome = "Status")

xgb_rec <- prep(xgb_rec, training = credit_data_tr)

bake(xgb_rec, credit_data_te, Assets)

```

step_embed*Encoding Factors into Multiple Columns*

Description

`step_embed()` creates a *specification* of a recipe step that will convert a nominal (i.e. factor) predictor into a set of scores derived from a tensorflow model via a word-embedding model. `embed_control` is a simple wrapper for setting default options.

Usage

```

step_embed(
  recipe,
  ...,
  role = "predictor",
  trained = FALSE,
  outcome = NULL,
  predictors = NULL,
  num_terms = 2,
  hidden_units = 0,
  options = embed_control(),
  mapping = NULL,
  history = NULL,
  keep_original_cols = FALSE,
  skip = FALSE,
  id = rand_id("embed")
)

embed_control(
  loss = "mse",
  metrics = NULL,
  optimizer = "sgd",
  epochs = 20,
  validation_split = 0,
  batch_size = 32,
  verbose = 0,
  callbacks = NULL
)

```

Arguments

recipe	A recipe object. The step will be added to the sequence of operations for this recipe.
...	One or more selector functions to choose variables. For <code>step_embed</code> , this indicates the variables to be encoded into a numeric format. See recipes::selections() for more details. For the <code>tidy</code> method, these are not currently used.
role	For model terms created by this step, what analysis role should they be assigned?. By default, the function assumes that the embedding variables created will be used as predictors in a model.
trained	A logical to indicate if the quantities for preprocessing have been estimated.
outcome	A call to <code>vars</code> to specify which variable is used as the outcome in the neural network.
predictors	An optional call to <code>vars</code> to specify any variables to be added as additional predictors in the neural network. These variables should be numeric and perhaps centered and scaled.
num_terms	An integer for the number of resulting variables.
hidden_units	An integer for the number of hidden units in a dense ReLu layer between the embedding and output later. Use a value of zero for no intermediate layer (see Details below).
options	A list of options for the model fitting process.
mapping	A list of tibble results that define the encoding. This is <code>NULL</code> until the step is trained by recipes::prep() .
history	A tibble with the convergence statistics for each term. This is <code>NULL</code> until the step is trained by recipes::prep() .
keep_original_cols	A logical to keep the original variables in the output. Defaults to <code>FALSE</code> .
skip	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations.
id	A character string that is unique to this step to identify it.
optimizer, loss, metrics	Arguments to pass to <code>keras3::compile()</code>
epochs, validation_split, batch_size, verbose, callbacks	Arguments to pass to <code>keras3::fit()</code>

Details

Factor levels are initially assigned at random to the new variables and these variables are used in a neural network to optimize both the allocation of levels to new columns as well as estimating a model to predict the outcome. See Section 6.1.2 of Francois and Allaire (2018) for more details.

The new variables are mapped to the specific levels seen at the time of model training and an extra instance of the variables are used for new levels of the factor.

One model is created for each call to `step_embed`. All terms given to the step are estimated and encoded in the same model which would also contain predictors given in `predictors` (if any).

When the outcome is numeric, a linear activation function is used in the last layer while softmax is used for factor outcomes (with any number of levels).

For example, the `keras3` code for a numeric outcome, one categorical predictor, and no hidden units used here would be

```
keras_model_sequential() |>
  layer_embedding(
    input_dim = num_factor_levels_x + 1,
    output_dim = num_terms
  ) |>
  layer_flatten() |>
  layer_dense(units = 1, activation = 'linear')
```

If a factor outcome is used and hidden units were requested, the code would be

```
keras_model_sequential() |>
  layer_embedding(
    input_dim = num_factor_levels_x + 1,
    output_dim = num_terms
  ) |>
  layer_flatten() |>
  layer_dense(units = hidden_units, activation = "relu") |>
  layer_dense(units = num_factor_levels_y, activation = 'softmax')
```

Other variables specified by `predictors` are added as an additional dense layer after `layer_flatten` and before the hidden layer.

Also note that it may be difficult to obtain reproducible results using this step due to the nature of Tensorflow (see link in References).

Tensorflow models cannot be run in parallel within the same session (via `foreach` or `futures`) or the `parallel` package. If using a recipes with this step with `caret`, avoid parallel processing.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables for encoding), `level` (the factor levels), and several columns containing `embed` in the name.

Tidying

When you `tidy()` this step, a tibble is returned with a number of columns with embedding information, and columns `terms`, `levels`, and `id`:

terms character, the selectors or variables selected

levels character, levels in variable

id character, id of this step

Tuning Parameters

This step has 2 tuning parameters:

- `num_terms`: # Model Terms (type: integer, default: 2)
- `hidden_units`: # Hidden Units (type: integer, default: 0)

Case weights

The underlying operation does not allow for case weights.

References

Francois C and Allaire JJ (2018) *Deep Learning with R*, Manning
 "Concatenate Embeddings for Categorical Variables with keras3" https://flovv.github.io/Embeddings_with_keras_part2/

Examples

```
data(grants, package = "modeldata")

set.seed(1)
grants_other <- sample_n(grants_other, 500)

rec <- recipe(class ~ num_ci + sponsor_code, data = grants_other) |>
  step_embed(sponsor_code,
    outcome = vars(class),
    options = embed_control(epochs = 10)
  )
```

step_lencode

Likelihood encoding using analytical formula

Description

`step_lencode()` creates a *specification* of a recipe step that will convert a nominal (i.e. factor) predictor into a single set of scores derived analytically.

Usage

```
step_lencode(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  smooth = TRUE,
  mapping = NULL,
```

```

  skip = FALSE,
  id = rand_id("lencode")
)

```

Arguments

recipe	A recipe object. The step will be added to the sequence of operations for this recipe.
...	One or more selector functions to choose variables. For <code>step_lencode()</code> , this indicates the variables to be encoded into a numeric format. See recipes::selections() for more details. For the <code>tidy</code> method, these are not currently used.
role	Not used by this step since no new variables are created.
trained	A logical to indicate if the quantities for preprocessing have been estimated.
outcome	A call to <code>vars</code> to specify which variable is used as the outcome. Only numeric and two-level factors are currently supported.
smooth	A logical, default to <code>TRUE</code> , should the estimates of groups with low counts be pulled towards the global estimate? Defaults to <code>TRUE</code> . See Details for how this is done. This is also known as partial pooling or shrinkage. Only works for numeric outcomes.
mapping	A list of tibble results that define the encoding. This is <code>NULL</code> until the step is trained by recipes::prep() .
skip	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

Each selected nominal predictor will be replaced by a numeric predictor. Each unique value of the nominal predictor is replaced by a numeric value. These values are calculated differently depending on the type of the outcome.

For **numeric** outcomes each value is the average value of the outcome inside each of the levels of the predictor. Unseen levels of the predictor will be using the global mean of the predictor. If case weights are used then a weighted mean is calculated instead.

For **nominal** outcomes each value is the log odds of the first level of the outcome variable being present, within each level of the levels of the predictor. Unseen levels will be replaced by the global log odds without stratification. If case weights are used then a weighted log odds is calculated.

If no or all occurrences happens then the log odds is calculated using $p = (2 * \text{nrow}(\text{data}) - 1) / (2 * \text{nrow}(\text{data}))$ to avoid infinity that would happen by taking the log of 0.

For numeric outcomes where `smooth = TRUE`, the following adjustment is done.

$$\text{estimate} = (n/\text{global_var}) / (n/\text{global_var} + 1/\text{outcome_var}) * \text{estimate} + (1/\text{outcome_var}) / (n/\text{global_var} + 1/\text{outcome_var}) * \text{global_mean}$$

Where n is the number of observations in the group.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables for encoding), `level` (the factor levels), and `value` (the encodings).

Tidying

When you `tidy()` this step, a tibble is returned with columns `level`, `value`, `terms`, and `id`:

level character, the factor levels
value numeric, the encoding
terms character, the selectors or variables selected
id character, id of this step

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in `recipes::case_weights` and the examples on tidymodels.org.

References

Micci-Barreca D (2001) "A preprocessing scheme for high-cardinality categorical attributes in classification and prediction problems," ACM SIGKDD Explorations Newsletter, 3(1), 27-32.
Zumel N and Mount J (2017) "vtreat: a data.frame Processor for Predictive Modeling," arXiv:1611.09477

See Also

`step_lencode_bayes()`, `step_lencode_glm()`, `step_lencode_mixed()`

Examples

```
library(recipes)
library(dplyr)
library(modeldata)

data(grants)

set.seed(1)
grants_other <- sample_n(grants_other, 500)
reencoded <- recipe(class ~ sponsor_code, data = grants_other) |>
  step_lencode(sponsor_code, outcome = vars(class), smooth = FALSE) |>
  prep()

bake(reencoded, grants_other)

tidy(reencoded, 1)
```

step_lencode_bayes*Supervised Factor Conversions into Linear Functions using Bayesian Likelihood Encodings*

Description

`step_lencode_bayes()` creates a *specification* of a recipe step that will convert a nominal (i.e. factor) predictor into a single set of scores derived from a generalized linear model estimated using Bayesian analysis.

Usage

```
step_lencode_bayes(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  options = list(seed = sample.int(10^5, 1)),
  verbose = FALSE,
  mapping = NULL,
  skip = FALSE,
  id = rand_id("lencode_bayes")
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose variables. For <code>step_lencode_bayes</code> , this indicates the variables to be encoded into a numeric format. See recipes::selections() for more details. For the <code>tidy</code> method, these are not currently used.
<code>role</code>	Not used by this step since no new variables are created.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>outcome</code>	A call to <code>vars</code> to specify which variable is used as the outcome in the generalized linear model. Only numeric and two-level factors are currently supported.
<code>options</code>	A list of options to pass to rstanarm::stan_glmer() .
<code>verbose</code>	A logical to control the default printing by rstanarm::stan_glmer() .
<code>mapping</code>	A list of tibble results that define the encoding. This is <code>NULL</code> until the step is trained by recipes::prep() .
<code>skip</code>	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
<code>id</code>	A character string that is unique to this step to identify it.

Details

For each factor predictor, a generalized linear model is fit to the outcome and the coefficients are returned as the encoding. These coefficients are on the linear predictor scale so, for factor outcomes, they are in log-odds units. The coefficients are created using a no intercept model and, when two factor outcomes are used, the log-odds reflect the event of interest being the *first* level of the factor.

For novel levels, a slightly trimmed average of the coefficients is returned.

A hierarchical generalized linear model is fit using `rstanarm::stan_glmer()` and no intercept via

```
stan_glmer(outcome ~ (1 | predictor), data = data, ...)
```

where the `...` include the `family` argument (automatically set by the step, unless passed in by options) as well as any arguments given to the `options` argument to the step. Relevant options include `chains`, `iter`, `cores`, and arguments for the priors (see the links in the References below). `prior_intercept` is the argument that has the most effect on the amount of shrinkage.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables for encoding), `level` (the factor levels), and `value` (the encodings).

Tidying

When you `tidy()` this step, a tibble is returned with columns `level`, `value`, `terms`, and `id`:

level character, the factor levels
value numeric, the encoding
terms character, the selectors or variables selected
id character, id of this step

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in `recipes::case_weights` and the examples on tidymodels.org.

References

- Micci-Barreca D (2001) "A preprocessing scheme for high-cardinality categorical attributes in classification and prediction problems," ACM SIGKDD Explorations Newsletter, 3(1), 27-32.
- Zumel N and Mount J (2017) "vtreat: a data.frame Processor for Predictive Modeling," arXiv:1611.09477 "Hierarchical Partial Pooling for Repeated Binary Trials" <https://CRAN.R-project.org/package=rstanarm/vignettes/pooling.html>
- "Prior Distributions for rstanarm Models" <http://mc-stan.org/rstanarm/reference/priors.html>
- "Estimating Generalized (Non-)Linear Models with Group-Specific Terms with rstanarm" <http://mc-stan.org/rstanarm/articles/glmer.html>

See Also

[step_lencode\(\)](#), [step_lencode_glm\(\)](#), [step_lencode_mixed\(\)](#)

Examples

```
library(recipes)
library(dplyr)
library(modeldata)

data(grants)

set.seed(1)
grants_other <- sample_n(grants_other, 500)

reencoded <- recipe(class ~ sponsor_code, data = grants_other) |>
  step_lencode_bayes(sponsor_code, outcome = vars(class))
```

step_lencode_glm

Supervised Factor Conversions into Linear Functions using Likelihood Encodings

Description

`step_lencode_glm()` creates a *specification* of a recipe step that will convert a nominal (i.e. factor) predictor into a single set of scores derived from a generalized linear model.

Usage

```
step_lencode_glm(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  mapping = NULL,
  skip = FALSE,
  id = rand_id("lencode_glm")
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose variables. For <code>step_lencode_glm</code> , this indicates the variables to be encoded into a numeric format. See recipes::selections() for more details. For the <code>tidy</code> method, these are not currently used.

role	Not used by this step since no new variables are created.
trained	A logical to indicate if the quantities for preprocessing have been estimated.
outcome	A call to <code>vars</code> to specify which variable is used as the outcome in the generalized linear model. Only numeric and two-level factors are currently supported.
mapping	A list of tibble results that define the encoding. This is <code>NULL</code> until the step is trained by <code>recipes::prep()</code> .
skip	A logical. Should the step be skipped when the recipe is baked by <code>recipes::bake()</code> ? While all operations are baked when <code>recipes::prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

For each factor predictor, a generalized linear model is fit to the outcome and the coefficients are returned as the encoding. These coefficients are on the linear predictor scale so, for factor outcomes, they are in log-odds units. The coefficients are created using a no intercept model and, when two factor outcomes are used, the log-odds reflect the event of interest being the *first* level of the factor.

For novel levels, a slightly trimmed average of the coefficients is returned.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables for encoding), `level` (the factor levels), and `value` (the encodings).

Tidying

When you `tidy()` this step, a tibble is returned with columns `level`, `value`, `terms`, and `id`:

- level** character, the factor levels
- value** numeric, the encoding
- terms** character, the selectors or variables selected
- id** character, id of this step

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in `recipes::case_weights` and the examples on tidymodels.org.

References

- Micci-Barreca D (2001) "A preprocessing scheme for high-cardinality categorical attributes in classification and prediction problems," *ACM SIGKDD Explorations Newsletter*, 3(1), 27-32.
- Zumel N and Mount J (2017) "vtreat: a data.frame Processor for Predictive Modeling," [arXiv:1611.09477](https://arxiv.org/abs/1611.09477)

See Also

[step_lencode\(\)](#), [step_lencode_bayes\(\)](#), [step_lencode_mixed\(\)](#)

Examples

```
library(recipes)
library(dplyr)
library(modeldata)

data(grants)

set.seed(1)
grants_other <- sample_n(grants_other, 500)

reencoded <- recipe(class ~ sponsor_code, data = grants_other) |>
  step_lencode_glm(sponsor_code, outcome = vars(class))
```

step_lencode_mixed	<i>Supervised Factor Conversions into Linear Functions using Bayesian Likelihood Encodings</i>
---------------------------	--

Description

`step_lencode_mixed()` creates a *specification* of a recipe step that will convert a nominal (i.e. factor) predictor into a single set of scores derived from a generalized linear mixed model.

Usage

```
step_lencode_mixed(
  recipe,
  ...,
  role = NA,
  trained = FALSE,
  outcome = NULL,
  options = list(verbose = 0),
  mapping = NULL,
  skip = FALSE,
  id = rand_id("lencode_mixed")
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
---------------------	--

...	One or more selector functions to choose variables. For <code>step_lencode_mixed</code> , this indicates the variables to be encoded into a numeric format. See <code>recipes::selections()</code> for more details. For the <code>tidy</code> method, these are not currently used.
<code>role</code>	Not used by this step since no new variables are created.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>outcome</code>	A call to <code>vars</code> to specify which variable is used as the outcome in the generalized linear model. Only numeric and two-level factors are currently supported.
<code>options</code>	A list of options to pass to <code>lme4::lmer()</code> or <code>lme4::glmer()</code> .
<code>mapping</code>	A list of tibble results that define the encoding. This is <code>NULL</code> until the step is trained by <code>recipes::prep()</code> .
<code>skip</code>	A logical. Should the step be skipped when the recipe is baked by <code>recipes::bake()</code> ? While all operations are baked when <code>recipes::prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
<code>id</code>	A character string that is unique to this step to identify it.

Details

For each factor predictor, a generalized linear model is fit to the outcome and the coefficients are returned as the encoding. These coefficients are on the linear predictor scale so, for factor outcomes, they are in log-odds units. The coefficients are created using a no intercept model and, when two factor outcomes are used, the log-odds reflect the event of interest being the *first* level of the factor.

For novel levels, a slightly timmed average of the coefficients is returned.

A hierarchical generalized linear model is fit using `lme4::lmer()` or `lme4::glmer()`, depending on the nature of the outcome, and no intercept via

```
lmer(outcome ~ 1 + (1 | predictor), data = data, ...)
```

where the `...` include the `family` argument (automatically set by the step) as well as any arguments given to the `options` argument to the step. Relevant options include `control` and `others`.

Value

An updated version of recipe with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables for encoding), `level` (the factor levels), and `value` (the encodings).

Tidying

When you `tidy()` this step, a tibble is returned with columns `level`, `value`, `terms`, and `id`:

level character, the factor levels

value numeric, the encoding

terms character, the selectors or variables selected

id character, id of this step

Case weights

This step performs an supervised operation that can utilize case weights. To use them, see the documentation in [recipes::case_weights](#) and the examples on [tidymodels.org](#).

References

Micci-Barreca D (2001) "A preprocessing scheme for high-cardinality categorical attributes in classification and prediction problems," ACM SIGKDD Explorations Newsletter, 3(1), 27-32.

Zumel N and Mount J (2017) "vtreat: a data.frame Processor for Predictive Modeling," arXiv:1611.09477

See Also

[step_lencode\(\)](#), [step_lencode_bayes\(\)](#), [step_lencode_glm\(\)](#)

Examples

```
library(recipes)
library(dplyr)
library(modeldata)

data(grants)

set.seed(1)
grants_other <- sample_n(grants_other, 500)

reencoded <- recipe(class ~ sponsor_code, data = grants_other) |>
  step_lencode_mixed(sponsor_code, outcome = vars(class))
```

step_pca_sparse *Sparse PCA Signal Extraction*

Description

`step_pca_sparse()` creates a *specification* of a recipe step that will convert numeric data into one or more principal components that can have some zero coefficients.

Usage

```
step_pca_sparse(
  recipe,
  ...,
  role = "predictor",
  trained = FALSE,
  num_comp = 5,
  predictor_prop = 1,
  options = list(),
```

```

  res = NULL,
  prefix = "PC",
  keep_original_cols = FALSE,
  skip = FALSE,
  id = rand_id("pca_sparse")
)

```

Arguments

recipe	A recipe object. The step will be added to the sequence of operations for this recipe.
...	One or more selector functions to choose which variables will be used to compute the components. See recipes::selections for more details. For the <code>tidy</code> method, these are not currently used.
role	For model terms created by this step, what analysis role should they be assigned? By default, the function assumes that the new principal component columns created by the original variables will be used as predictors in a model.
trained	A logical to indicate if the quantities for preprocessing have been estimated.
num_comp	The number of components to retain as new predictors. If <code>num_comp</code> is greater than the number of columns or the number of possible components, a smaller value will be used. If <code>num_comp = 0</code> is set then no transformation is done and selected variables will stay unchanged, regardless of the value of <code>keep_original_cols</code> .
predictor_prop	The maximum number of original predictors that can have non-zero coefficients for each PCA component (via regularization).
options	A list of options to the default method for irlba::ssvd() .
res	The rotation matrix once this preprocessing step has been trained by recipes::prep .
prefix	A character string that will be the prefix to the resulting new variables. See notes below.
keep_original_cols	A logical to keep the original variables in the output. Defaults to FALSE.
skip	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

The `irlba` package is required for this step. If it is not installed, the user will be prompted to do so when the step is defined. The `irlba::ssvd()` function is used to encourage sparsity; that documentation has details about this method.

The argument `num_comp` controls the number of components that will be retained (the original variables that are used to derive the components are removed from the data). The new components will have names that begin with `prefix` and a sequence of numbers. The variable names are padded with zeros. For example, if `num_comp < 10`, their names will be PC1 - PC9. If `num_comp = 101`, the names would be PC1 - PC101.

Value

An updated version of `recipe` with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables selected), `value` (the loading), and `component`.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `value`, `component`, and `id`:

terms character, the selectors or variables selected
value numeric, variable loading
component character, principle component
id character, id of this step

Tuning Parameters

This step has 2 tuning parameters:

- `num_comp`: # Components (type: integer, default: 5)
- `predictor_prop`: Proportion of Predictors (type: double, default: 1)

Case weights

The underlying operation does not allow for case weights.

See Also

[recipes::step_pca\(\)](#), [recipes::step_kpca\(\)](#), [recipes::step_kpca_poly\(\)](#), [recipes::step_kpca_rbf\(\)](#), [step_pca_sparse_bayes\(\)](#), [step_pca_truncated\(\)](#)

Examples

```
library(recipes)
library(ggplot2)

data(ad_data, package = "modeldata")

ad_rec <-
  recipe(Class ~ ., data = ad_data) |>
  step_zv(all_predictors()) |>
  step_YeoJohnson(all_numeric_predictors()) |>
  step_normalize(all_numeric_predictors()) |>
  step_pca_sparse(
    all_numeric_predictors(),
    predictor_prop = 0.75,
    num_comp = 3,
    id = "sparse pca"
  ) |>
  prep()
```

```

tidy(ad_rec, id = "sparse pca") |>
  mutate(value = ifelse(value == 0, NA, value)) |>
  ggplot(aes(x = component, y = terms, fill = value)) +
  geom_tile() +
  scale_fill_gradient2() +
  theme(axis.text.y = element_blank())

```

step_pca_sparse_bayes *Sparse Bayesian PCA Signal Extraction*

Description

`step_pca_sparse_bayes()` creates a *specification* of a recipe step that will convert numeric data into one or more principal components that can have some zero coefficients.

Usage

```

step_pca_sparse_bayes(
  recipe,
  ...,
  role = "predictor",
  trained = FALSE,
  num_comp = 5,
  prior_slab_dispersion = 1,
  prior_mixture_threshold = 0.1,
  options = list(),
  res = NULL,
  prefix = "PC",
  keep_original_cols = FALSE,
  skip = FALSE,
  id = rand_id("pca_sparse_bayes")
)

```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose which variables will be used to compute the components. See recipes::selections for more details. For the <code>tidy</code> method, these are not currently used.
<code>role</code>	For model terms created by this step, what analysis role should they be assigned? By default, the function assumes that the new principal component columns created by the original variables will be used as predictors in a model.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.

num_comp	The number of components to retain as new predictors. If num_comp is greater than the number of columns or the number of possible components, a smaller value will be used. If num_comp = 0 is set then no transformation is done and selected variables will stay unchanged, regardless of the value of keep_original_cols.
prior_slab_dispersion	This value is proportional to the dispersion (or scale) parameter for the slab portion of the prior. Smaller values result in an increase in zero coefficients.
prior_mixture_threshold	The parameter that defines the trade-off between the spike and slab components of the prior. Increasing this parameter increases the number of zero coefficients.
options	A list of options to the default method for <code>VBsparsePCA::VBsparsePCA()</code> .
res	The rotation matrix once this preprocessing step has been trained by <code>recipes::prep()</code> .
prefix	A character string that will be the prefix to the resulting new variables. See notes below.
keep_original_cols	A logical to keep the original variables in the output. Defaults to FALSE.
skip	A logical. Should the step be skipped when the recipe is baked by <code>recipes::bake()</code> ? While all operations are baked when <code>recipes::prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using skip = TRUE as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

The `VBsparsePCA` package is required for this step. If it is not installed, the user will be prompted to do so when the step is defined.

A spike-and-slab prior is a mixture of two priors. One (the "spike") has all of its mass at zero and represents a variable that has no contribution to the PCA coefficients. The other prior is a broader distribution that reflects the coefficient distribution of variables that do affect the PCA analysis. This is the "slab". The narrower the slab, the more likely that a coefficient will be zero (or are regularized to be closer to zero). The mixture of these two priors is governed by a mixing parameter, which itself has a prior distribution and a hyper-parameter prior.

PCA coefficients and their resulting scores are unique only up to the sign. This step will attempt to make the sign of the components more consistent from run-to-run. However, the sparsity constraint may interfere with this goal.

The argument num_comp controls the number of components that will be retained (the original variables that are used to derive the components are removed from the data). The new components will have names that begin with prefix and a sequence of numbers. The variable names are padded with zeros. For example, if num_comp < 10, their names will be PC1 - PC9. If num_comp = 101, the names would be PC1 - PC101.

Value

An updated version of recipe with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with columns `terms` (the selectors or variables selected), `value` (the loading), and `component`.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms`, `value`, `component`, and `id`:

terms character, the selectors or variables selected
value numeric, variable loading
component character, principle component
id character, id of this step

Tuning Parameters

This step has 3 tuning parameters:

- `num_comp`: # Components (type: integer, default: 5)
- `prior_slab_dispersion`: Dispersion of Slab Prior (type: double, default: 1)
- `prior_mixture_threshold`: Threshold for Mixture Prior (type: double, default: 0.1)

Case weights

The underlying operation does not allow for case weights.

References

Ning, B. (2021). Spike and slab Bayesian sparse principal component analysis. arXiv:2102.00305.

See Also

`recipes::step_pca()`, `recipes::step_kpca()`, `recipes::step_kpca_poly()`, `recipes::step_kpca_rbf()`, `step_pca_sparse()`, `step_pca_truncated()`

Examples

```
library(recipes)
library(ggplot2)

data(ad_data, package = "modeldata")

ad_rec <-
  recipe(Class ~ ., data = ad_data) |>
  step_zv(all_predictors()) |>
  step_YeoJohnson(all_numeric_predictors()) |>
  step_normalize(all_numeric_predictors()) |>
  step_pca_sparse_bayes(
    all_numeric_predictors(),
    prior_mixture_threshold = 0.95,
    prior_slab_dispersion = 0.05,
    num_comp = 3,
    id = "sparse bayesian pca"
  ) |>
  prep()
```

```
tidy(ad_rec, id = "sparse bayesian pca") |>
  mutate(value = ifelse(value == 0, NA, value)) |>
  ggplot(aes(x = component, y = terms, fill = value)) +
  geom_tile() +
  scale_fill_gradient2() +
  theme(axis.text.y = element_blank())
```

step_pca_truncated *Truncated PCA Signal Extraction*

Description

`step_pca_truncated()` creates a *specification* of a recipe step that will convert numeric data into one or more principal components. It is truncated as it only calculates the number of components it is asked instead of all of them as is done in `recipes::step_pca()`.

Usage

```
step_pca_truncated(
  recipe,
  ...,
  role = "predictor",
  trained = FALSE,
  num_comp = 5,
  options = list(),
  res = NULL,
  columns = NULL,
  prefix = "PC",
  keep_original_cols = FALSE,
  skip = FALSE,
  id = rand_id("pca_truncated")
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose variables for this step. See <code>selections()</code> for more details.
<code>role</code>	For model terms created by this step, what analysis role should they be assigned? By default, the new columns created by this step from the original variables will be used as <i>predictors</i> in a model.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.
<code>num_comp</code>	The number of components to retain as new predictors. If <code>num_comp</code> is greater than the number of columns or the number of possible components, a smaller value will be used. If <code>num_comp = 0</code> is set then no transformation is done and selected variables will stay unchanged, regardless of the value of <code>keep_original_cols</code> .

options	A list of options to the default method for <code>irlba::prcomp_irlba()</code> . Argument defaults are set to <code>retx = FALSE</code> , <code>center = FALSE</code> , <code>scale. = FALSE</code> , and <code>tol = NULL</code> . Note that the argument <code>x</code> should not be passed here (or at all).
res	The <code>irlba::prcomp_irlba()</code> object is stored here once this preprocessing step has been trained by <code>recipes::prep</code> .
columns	A character string of the selected variable names. This field is a placeholder and will be populated once <code>prep()</code> is used.
prefix	A character string for the prefix of the resulting new variables. See notes below.
keep_original_cols	A logical to keep the original variables in the output. Defaults to <code>FALSE</code> .
skip	A logical. Should the step be skipped when the recipe is baked by <code>bake()</code> ? While all operations are baked when <code>prep()</code> is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations.
id	A character string that is unique to this step to identify it.

Details

Principal component analysis (PCA) is a transformation of a group of variables that produces a new set of artificial features or components. These components are designed to capture the maximum amount of information (i.e. variance) in the original variables. Also, the components are statistically independent from one another. This means that they can be used to combat large inter-variables correlations in a data set.

It is advisable to standardize the variables prior to running PCA. Here, each variable will be centered and scaled prior to the PCA calculation. This can be changed using the `options` argument or by using `recipes::step_center()` and `recipes::step_scale()`.

The argument `num_comp` controls the number of components that will be retained (the original variables that are used to derive the components are removed from the data). The new components will have names that begin with `prefix` and a sequence of numbers. The variable names are padded with zeros. For example, if `num_comp < 10`, their names will be `PC1 - PC9`. If `num_comp = 101`, the names would be `PC1 - PC101`.

Value

An updated version of `recipe` with the new step added to the sequence of any existing operations.

Tidying

When you `tidy()` this step two things can happen depending the `type` argument. If `type = "coef"` a tibble returned with 4 columns `terms`, `value`, `component`, and `id`:

- terms** character, the selectors or variables selected
- value** numeric, variable loading
- component** character, principle component
- id** character, id of this step

If `type = "variance"` a tibble returned with 4 columns `terms`, `value`, `component`, and `id`:

terms character, type of variance
value numeric, value of the variance
component integer, principle component
id character, id of this step

Tuning Parameters

This step has 1 tuning parameters:

- `num_comp`: # Components (type: integer, default: 5)

Case weights

This step performs an unsupervised operation that can utilize case weights. As a result, case weights are only used with frequency weights. For more information, see the documentation in [recipes::case_weights](#) and the examples on [tidymodels.org](#).

References

Jolliffe, I. T. (2010). *Principal Component Analysis*. Springer.

See Also

[recipes::step_pca\(\)](#), [recipes::step_kpca\(\)](#), [recipes::step_kpca_poly\(\)](#), [recipes::step_kpca_rbf\(\)](#), [step_pca_sparse_bayes\(\)](#), [step_pca_sparse\(\)](#)

Examples

```
rec <- recipe(~., data = mtcars)
pca_trans <- rec |>
  step_normalize(all_numeric()) |>
  step_pca_truncated(all_numeric(), num_comp = 2)
pca_estimates <- prep(pca_trans, training = mtcars)
pca_data <- bake(pca_estimates, mtcars)

rng <- extendrange(c(pca_data$PC1, pca_data$PC2))
plot(pca_data$PC1, pca_data$PC2,
  xlim = rng, ylim = rng
)

tidy(pca_trans, number = 2)
tidy(pca_estimates, number = 2)
```

step_umap	<i>Supervised and unsupervised uniform manifold approximation and projection (UMAP)</i>
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Description

`step_umap()` creates a *specification* of a recipe step that will project a set of features into a smaller space.

Usage

```
step_umap(  
  recipe,  
  ...,  
  role = "predictor",  
  trained = FALSE,  
  outcome = NULL,  
  neighbors = 15,  
  num_comp = 2,  
  min_dist = 0.01,  
  metric = "euclidean",  
  learn_rate = 1,  
  epochs = NULL,  
  initial = "spectral",  
  target_weight = 0.5,  
  options = list(verbose = FALSE, n_threads = 1),  
  seed = sample(10^5, 2),  
  prefix = "UMAP",  
  keep_original_cols = FALSE,  
  retain = deprecated(),  
  object = NULL,  
  skip = FALSE,  
  id = rand_id("umap")  
)
```

Arguments

<code>recipe</code>	A recipe object. The step will be added to the sequence of operations for this recipe.
<code>...</code>	One or more selector functions to choose variables for this step. See selections() for more details.
<code>role</code>	For model terms created by this step, what analysis role should they be assigned? By default, the new columns created by this step from the original variables will be used as <i>predictors</i> in a model.
<code>trained</code>	A logical to indicate if the quantities for preprocessing have been estimated.

outcome	A call to <code>vars</code> to specify which variable is used as the outcome in the encoding process (if any).
neighbors	An integer for the number of nearest neighbors used to construct the target simplicial set. If <code>neighbors</code> is greater than the number of data points, the smaller value is used.
num_comp	An integer for the number of UMAP components. If <code>num_comp</code> is greater than the number of selected columns minus one, the smaller value is used.
min_dist	The effective minimum distance between embedded points.
metric	Character, type of distance metric to use to find nearest neighbors. See uwot::umap() for more details. Default to "euclidean".
learn_rate	Positive number of the learning rate for the optimization process.
epochs	Number of iterations for the neighbor optimization. See uwot::umap() for more details.
initial	Character, Type of initialization for the coordinates. Can be one of "spectral", "normlaplacian", "random", "lvrandom", "laplacian", "pca", "spca", "agspectral", or a matrix of initial coordinates. See uwot::umap() for more details. Default to "spectral".
target_weight	Weighting factor between data topology and target topology. A value of 0.0 weights entirely on data, a value of 1.0 weights entirely on target. The default of 0.5 balances the weighting equally between data and target.
options	A list of options to pass to uwot::umap() . The arguments <code>X</code> , <code>n_neighbors</code> , <code>n_components</code> , <code>min_dist</code> , <code>n_epochs</code> , <code>ret_model</code> , and <code>learning_rate</code> should not be passed here. By default, <code>verbose</code> and <code>n_threads</code> are set.
seed	Two integers to control the random numbers used by the numerical methods. The default pulls from the main session's stream of numbers and will give reproducible results if the seed is set prior to calling recipes::prep or recipes::bake .
prefix	A character string for the prefix of the resulting new variables. See notes below.
keep_original_cols	A logical to keep the original variables in the output. Defaults to FALSE.
retain	Use <code>keep_original_cols</code> instead to specify whether the original predictors should be retained along with the new embedding variables.
object	An object that defines the encoding. This is <code>NULL</code> until the step is trained by recipes::prep() .
skip	A logical. Should the step be skipped when the recipe is baked by bake() ? While all operations are baked when prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations.
id	A character string that is unique to this step to identify it.

Details

UMAP, short for Uniform Manifold Approximation and Projection, is a nonlinear dimension reduction technique that finds local, low-dimensional representations of the data. It can be run unsupervised or supervised with different types of outcome data (e.g. numeric, factor, etc).

The argument `num_comp` controls the number of components that will be retained (the original variables that are used to derive the components are removed from the data). The new components will have names that begin with `prefix` and a sequence of numbers. The variable names are padded with zeros. For example, if `num_comp < 10`, their names will be UMAP1 - UMAP9. If `num_comp = 101`, the names would be UMAP1 - UMAP101.

Value

An updated version of `recipe` with the new step added to the sequence of any existing operations.

Tidying

When you `tidy()` this step, a tibble is returned with columns `terms` and `id`:

`terms` character, the selectors or variables selected

`id` character, id of this step

Tuning Parameters

This step has 7 tuning parameters:

- `num_comp`: # Components (type: integer, default: 2)
- `neighbors`: # Nearest Neighbors (type: integer, default: 15)
- `min_dist`: Min Distance between Points (type: double, default: 0.01)
- `learn_rate`: Learning Rate (type: double, default: 1)
- `epochs`: # Epochs (type: integer, default: NULL)
- `initial`: UMAP Initialization (type: character, default: spectral)
- `target_weight`: Proportion Supervised (type: double, default: 0.5)

Case weights

The underlying operation does not allow for case weights.

Saving prepped recipe object

This recipe step may require native serialization when saving for use in another R session. To learn more about serialization for prepped recipes, see the `bundle` package.

References

McInnes, L., & Healy, J. (2018). UMAP: Uniform Manifold Approximation and Projection for Dimension Reduction. <https://arxiv.org/abs/1802.03426>.

"How UMAP Works" https://umap-learn.readthedocs.io/en/latest/how_umap_works.html

Examples

```

library(recipes)
library(ggplot2)

split <- seq.int(1, 150, by = 9)
tr <- iris[-split, ]
te <- iris[split, ]

set.seed(11)
supervised <-
  recipe(Species ~ ., data = tr) |>
  step_center(all_predictors()) |>
  step_scale(all_predictors()) |>
  step_umap(all_predictors(), outcome = vars(Species), num_comp = 2) |>
  prep(training = tr)

theme_set(theme_bw())

bake(supervised, new_data = te, Species, starts_with("umap")) |>
  ggplot(aes(x = UMAP1, y = UMAP2, col = Species)) +
  geom_point(alpha = .5)

```

step_woe

Weight of evidence transformation

Description

`step_woe()` creates a *specification* of a recipe step that will transform nominal data into its numerical transformation based on weights of evidence against a binary outcome.

Usage

```

step_woe(
  recipe,
  ...,
  role = "predictor",
  outcome,
  trained = FALSE,
  dictionary = NULL,
  Laplace = 1e-06,
  prefix = "woe",
  keep_original_cols = FALSE,
  skip = FALSE,
  id = rand_id("woe")
)

```

Arguments

recipe	A recipe object. The step will be added to the sequence of operations for this recipe.
...	One or more selector functions to choose which variables will be used to compute the components. See recipes::selections for more details. For the <code>tidy</code> method, these are not currently used.
role	For model terms created by this step, what analysis role should they be assigned?. By default, the function assumes that the new woe components columns created by the original variables will be used as predictors in a model.
outcome	The bare name of the binary outcome encased in <code>vars()</code> .
trained	A logical to indicate if the quantities for preprocessing have been estimated.
dictionary	A <code>tbl</code> . A map of levels and woe values. It must have the same layout than the output returned from dictionary() . If <code>NULL</code> the function will build a dictionary with those variables passed to See dictionary() for details.
Laplace	The Laplace smoothing parameter. A value usually applied to avoid -Inf/Inf from predictor category with only one outcome class. Set to 0 to allow Inf/Inf. The default is <code>1e-6</code> . Also known as 'pseudocount' parameter of the Laplace smoothing technique.
prefix	A character string that will be the prefix to the resulting new variables. See notes below.
keep_original_cols	A logical to keep the original variables in the output. Defaults to <code>FALSE</code> .
skip	A logical. Should the step be skipped when the recipe is baked by recipes::bake() ? While all operations are baked when recipes::prep() is run, some operations may not be able to be conducted on new data (e.g. processing the outcome variable(s)). Care should be taken when using <code>skip = TRUE</code> as it may affect the computations for subsequent operations
id	A character string that is unique to this step to identify it.

Details

WoE is a transformation of a group of variables that produces a new set of features. The formula is

$$woe_c = \log((P(X = c|Y = 1))/(P(X = c|Y = 0)))$$

where c goes from 1 to C levels of a given nominal predictor variable X .

These components are designed to transform nominal variables into numerical ones with the property that the order and magnitude reflects the association with a binary outcome. To apply it on numerical predictors, it is advisable to discretize the variables prior to running WoE. Here, each variable will be binarized to have woe associated later. This can be achieved by using [recipes::step_discretize\(\)](#).

The argument `Laplace` is a small quantity added to the proportions of 1's and 0's with the goal to avoid $\log(p/0)$ or $\log(0/p)$ results. The numerical woe versions will have names that begin with `woe_` followed by the respective original name of the variables. See Good (1985).

One can pass a custom `dictionary` tibble to `step_woe()`. It must have the same structure of the output from `dictionary()` (see examples). If not provided it will be created automatically. The

role of this tibble is to store the map between the levels of nominal predictor to its woe values. You may want to tweak this object with the goal to fix the orders between the levels of one given predictor. One easy way to do this is by tweaking an output returned from `dictionary()`.

Value

An updated version of recipe with the new step added to the sequence of existing steps (if any). For the `tidy` method, a tibble with the woe dictionary used to map categories with woe values.

Tidying

When you `tidy()` this step, a tibble with columns `terms` (the selectors or variables selected), `value`, `n_tot`, `n_bad`, `n_good`, `p_bad`, `p_good`, `woe` and `outcome` is returned.. See `dictionary()` for more information.

When you `tidy()` this step, a tibble is returned with columns `value`, `n_tot`, `n_bad`, `n_good`, `p_bad`, `p_good`, `woe` and `outcome` and `id`:

terms character, the selectors or variables selected

value character, level of the outcome

n_tot integer, total number

n_bad integer, number of bad examples

n_good integer, number of good examples

p_bad numeric, p of bad examples

p_good numeric, p of good examples

woe numeric, weight of evidence

outcome character, name of outcome variable

id character, id of this step

Tuning Parameters

This step has 1 tuning parameters:

- **Laplace**: Laplace Correction (type: double, default: 1e-06)

Case weights

The underlying operation does not allow for case weights.

References

Kullback, S. (1959). *Information Theory and Statistics*. Wiley, New York.

Hastie, T., Tibshirani, R. and Friedman, J. (1986). *Elements of Statistical Learning*, Second Edition, Springer, 2009.

Good, I. J. (1985), "Weight of evidence: A brief survey", *Bayesian Statistics*, 2, pp.249-270.

Examples

```

library(modeldata)
data("credit_data")

set.seed(111)
in_training <- sample(1:nrow(credit_data), 2000)

credit_tr <- credit_data[in_training, ]
credit_te <- credit_data[-in_training, ]

rec <- recipe(Status ~ ., data = credit_tr) |>
  step_woe(Job, Home, outcome = vars(Status))

woe_models <- prep(rec, training = credit_tr)

# the encoding:
bake(woe_models, new_data = credit_te |> slice(1:5), starts_with("woe"))
# the original data
credit_te |>
  slice(1:5) |>
  dplyr::select(Job, Home)
# the details:
tidy(woe_models, number = 1)

# Example of custom dictionary + tweaking
# custom dictionary
woe_dict_custom <- credit_tr |> dictionary(Job, Home, outcome = "Status")
woe_dict_custom[4, "woe"] <- 1.23 # tweak

# passing custom dict to step_woe()
rec_custom <- recipe(Status ~ ., data = credit_tr) |>
  step_woe(
    Job, Home,
    outcome = vars(Status), dictionary = woe_dict_custom
  ) |>
  prep()

rec_custom_baked <- bake(rec_custom, new_data = credit_te)
rec_custom_baked |>
  dplyr::filter(woe_Job == 1.23) |>
  head()

```

woe_table

Crosstable with woe between a binary outcome and a predictor variable.

Description

Calculates some summaries and the WoE (Weight of Evidence) between a binary outcome and a given predictor variable. Used to build the dictionary.

Usage

```
woe_table(predictor, outcome, Laplace = 1e-06, call = rlang::caller_env(0))
```

Arguments

<code>predictor</code>	A atomic vector, usually with few distinct values.
<code>outcome</code>	The dependent variable. A atomic vector with exactly 2 distinct values.
<code>Laplace</code>	The pseudocount parameter of the Laplace Smoothing estimator. Default to 1e-6. Value to avoid -Inf/Inf from predictor category with only one outcome class. Set to 0 to allow Inf/-Inf.
<code>call</code>	The execution environment of a currently running function, e.g. <code>caller_env()</code> . The function will be mentioned in error messages as the source of the error. See the <code>call</code> argument of rlang::abort() for more information.

Value

a tibble with counts, proportions and woe. Warning: woe can possibly be -Inf. Use 'Laplace' arg to avoid that.

References

Kullback, S. (1959). *Information Theory and Statistics*. Wiley, New York.

Hastie, T., Tibshirani, R. and Friedman, J. (1986). *Elements of Statistical Learning*, Second Edition, Springer, 2009.

Good, I. J. (1985), "Weight of evidence: A brief survey", *Bayesian Statistics*, 2, pp.249-270.

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