Package ‘hal9001’

November 14, 2023

Title  The Scalable Highly Adaptive Lasso
Version  0.4.6
Description  A scalable implementation of the highly adaptive lasso algorithm, including routines for constructing sparse matrices of basis functions of the observed data, as well as a custom implementation of Lasso regression tailored to enhance efficiency when the matrix of predictors is composed exclusively of indicator functions. For ease of use and increased flexibility, the Lasso fitting routines invoke code from the ‘glmnet’ package by default. The highly adaptive lasso was first formulated and described by MJ van der Laan (2017) <doi:10.1515/ijb-2015-0097>, with practical demonstrations of its performance given by Benkeser and van der Laan (2016) <doi:10.1109/DSAA.2016.93>. This implementation of the highly adaptive lasso algorithm was described by Hejazi, Coyle, and van der Laan (2020) <doi:10.21105/joss.02526>.

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HAL Formula addition: Adding formula term object together into a single formula object term.

Description

HAL Formula addition: Adding formula term object together into a single formula object term.

Usage

```r
## S3 method for class 'formula_hal9001'

x + y
```

Arguments

- `x`: A `formula_hal9001` object as outputted by `h`.
- `y`: A `formula_hal9001` object as outputted by `h`.

apply_copy_map

Apply copy map

Description

OR duplicate training set columns together

Usage

```r
apply_copy_map(X, copy_map)
```

Arguments

- `X`: Sparse matrix containing columns of indicator functions.
- `copy_map`: the copy map

Value

A `dgCMatrix` sparse matrix corresponding to the design matrix for a zero-th order highly adaptive lasso, but with all duplicated columns (basis functions) removed.
Examples

gendata <- function(n) {
  W1 <- runif(n, -3, 3)
  W2 <- rnorm(n)
  W3 <- runif(n)
  W4 <- rnorm(n)
  g0 <- plogis(0.5 * (-0.8 * W1 + 0.39 * W2 + 0.08 * W3 - 0.12 * W4))
  A <- rbinom(n, 1, g0)
  Q0 <- plogis(0.15 * (2 * A + 2 * A * W1 + 6 * A * W3 * W4 - 3))
  Y <- rbinom(n, 1, Q0)
  data.frame(A, W1, W2, W3, W4, Y)
}
set.seed(1234)
data <- gendata(100)
covars <- setdiff(names(data), "Y")
X <- as.matrix(data[, covars, drop = FALSE])
basis_list <- enumerate_basis(X)
x_basis <- make_design_matrix(X, basis_list)
copy_map <- make_copy_map(x_basis)
x_basis_uniq <- apply_copy_map(x_basis, copy_map)

as_dgCMatrix 

Fast Coercion to Sparse Matrix

Description


Usage

as_dgCMatrix(XX_)

Arguments

XX_ An object of class Matrix that has a sparse structure suitable for coercion to a sparse matrix format of dgCMatrix.

Value

An object of class dgCMatrix, coerced from input XX_.

basis_list_cols

List Basis Functions

Description
Build a list of basis functions from a set of columns

Usage

basis_list_cols(
cols,
  x,
  smoothness_orders,
  include_zero_order,
  include_lower_order = FALSE
)

Arguments

cols Index or indices (as numeric) of covariates (columns) of interest in the data matrix x for which basis functions ought to be generated. Note that basis functions for interactions of these columns are computed automatically.

x A matrix containing observations in the rows and covariates in the columns. Basis functions are computed for these covariates.

smoothness_orders An integer vector of length ncol(x) specifying the desired smoothness of the function in each covariate. k = 0 is no smoothness (indicator basis), k = 1 is first order smoothness, and so on. For an additive model, the component function for each covariate will have the degree of smoothness as specified by smoothness_orders. For non-additive components (tensor products of univariate basis functions), the univariate basis functions in each tensor product have smoothness degree as specified by smoothness_orders.

include_zero_order A logical, indicating whether the zeroth order basis functions are included for each covariate (if TRUE), in addition to the smooth basis functions given by smoothness_orders. This allows the algorithm to data-adaptively choose the appropriate degree of smoothness.

include_lower_order A logical, like include_zero_order, except including all basis functions of lower smoothness degrees than specified via smoothness_orders.

Value
A list containing the basis functions generated from a set of input columns.
basis_of_degree  

**Compute Degree of Basis Functions**

**Description**

Find the full list of basis functions up to a particular degree

**Usage**

```r
basis_of_degree(
  x,
  degree,
  smoothness_orders,
  include_zero_order,
  include_lower_order
)
```

**Arguments**

- `x`: An input matrix containing observations and covariates following standard conventions in problems of statistical learning.
- `degree`: The highest order of interaction terms for which the basis functions ought to be generated. The default (NULL) corresponds to generating basis functions for the full dimensionality of the input matrix.
- `smoothness_orders`: An integer vector of length `ncol(x)` specifying the desired smoothness of the function in each covariate. k = 0 is no smoothness (indicator basis), k = 1 is first order smoothness, and so on. For an additive model, the component function for each covariate will have the degree of smoothness as specified by smoothness_orders. For non-additive components (tensor products of univariate basis functions), the univariate basis functions in each tensor product have smoothness degree as specified by smoothness_orders.
- `include_zero_order`: A logical, indicating whether the zeroth order basis functions are included for each covariate (if TRUE), in addition to the smooth basis functions given by smoothness_orders. This allows the algorithm to data-adaptively choose the appropriate degree of smoothness.
- `include_lower_order`: A logical, like include_zero_order, except including all basis functions of lower smoothness degrees than specified via smoothness_orders.

**Value**

A list containing basis functions and cutoffs generated from a set of input columns up to a particular pre-specified degree.
enumerate_basis

Enumerate Basis Functions

Description
Generate basis functions for all covariates and interaction terms thereof up to a specified order/degree.

Usage
enumerate_basis(
  x,
  max_degree = NULL,
  smoothness_orders = rep(0, ncol(x)),
  include_zero_order = FALSE,
  include_lower_order = FALSE,
  num_knots = NULL
)

Arguments
x                An input matrix containing observations and covariates following standard conventions in problems of statistical learning.
max_degree       The highest order of interaction terms for which the basis functions ought to be generated. The default (NULL) corresponds to generating basis functions for the full dimensionality of the input matrix.
smoothness_orders An integer vector of length ncol(x) specifying the desired smoothness of the function in each covariate. k = 0 is no smoothness (indicator basis), k = 1 is first order smoothness, and so on. For an additive model, the component function for each covariate will have the degree of smoothness as specified by smoothness_orders. For non-additive components (tensor products of univariate basis functions), the univariate basis functions in each tensor product have smoothness degree as specified by smoothness_orders.
include_zero_order A logical, indicating whether the zeroth order basis functions are included for each covariate (if TRUE), in addition to the smooth basis functions given by smoothness_orders. This allows the algorithm to data-adaptively choose the appropriate degree of smoothness.
include_lower_order A logical, like include_zero_order, except including all basis functions of lower smoothness degrees than specified via smoothness_orders.
num_knots        A vector of length max_degree, which determines how granular the knot points to generate basis functions should be for each degree of basis function. The first entry of num_knots determines the number of knot points to be used for each univariate basis function. More generally, The kth entry of num_knots determines the number of knot points to be used for the kth degree basis functions.
Specifically, for a kth degree basis function, which is the tensor product of k univariate basis functions, this determines the number of knot points to be used for each univariate basis function in the tensor product.

Value

A list of basis functions generated for all covariates and interaction thereof up to a pre-specified degree.

Examples

gendata <- function(n) {
  W1 <- runif(n, -3, 3)
  W2 <- rnorm(n)
  W3 <- runif(n)
  W4 <- rnorm(n)
  g0 <- plogis(0.5 * (-0.8 * W1 + 0.39 * W2 + 0.08 * W3 - 0.12 * W4))
  A <- rbinom(n, 1, g0)
  Q0 <- plogis(0.15 * (2 * A + 2 * A * W1 + 6 * A * W3 * W4 - 3))
  Y <- rbinom(n, 1, Q0)
  data.frame(A, W1, W2, W3, W4, Y)
}
set.seed(1234)
data <- gendata(100)
covars <- setdiff(names(data), "Y")
X <- as.matrix(data[, covars, drop = FALSE])
basis_list <- enumerate_basis(X)

---

**evaluate_basis**

**Generate Basis Functions**

**Description**

Populates a column (indexed by basis_col) of x.basis with basis indicators.

**Usage**

evaluate_basis(basis, X, x.basis, basis_col)

**Arguments**

- **basis**: The basis function.
- **X**: The design matrix, containing the original data.
- **x.basis**: The HAL design matrix, containing indicator functions.
- **basis_col**: Numeric indicating which column to populate.
Description

Estimation procedure for HAL, the Highly Adaptive Lasso

Usage

```r
fit_hal(
  X,
  Y,
  formula = NULL,
  X_unpenalized = NULL,
  max_degree = ifelse(ncol(X) >= 20, 2, 3),
  smoothness_orders = 1,
  num_knots = num_knots_generator(max_degree = max_degree, smoothness_orders =
    smoothness_orders, base_num_knots_0 = 200, base_num_knots_1 = 50),
  reduce_basis = NULL,
  family = c("gaussian", "binomial", "poisson", "cox", "mgaussian"),
  lambda = NULL,
  id = NULL,
  weights = NULL,
  offset = NULL,
  fit_control = list(cv_select = TRUE, use_min = TRUE, lambda.min.ratio = 1e-04,
    prediction_bounds = "default"),
  basis_list = NULL,
  return_lasso = TRUE,
  return_xbasis = FALSE,
  yolo = FALSE
)
```

Arguments

- **X**: An input matrix with dimensions number of observations-by-number of covariates that will be used to derive the design matrix of basis functions.
- **Y**: A numeric vector of observations of the outcome variable. For family="mgaussian", Y is a matrix of observations of the outcome variables.
- **formula**: A character string formula to be used in `formula_hal`. See its documentation for details.
- **X_unpenalized**: An input matrix with the same number of rows as X, for which no L1 penalization will be performed. Note that X_unpenalized is directly appended to the design matrix; no basis expansion is performed on X_unpenalized.
- **max_degree**: The highest order of interaction terms for which basis functions ought to be generated.
smoothness_orders

An integer, specifying the smoothness of the basis functions. See details for smoothness_orders for more information.

num_knots

An integer vector of length 1 or max_degree, specifying the maximum number of knot points (i.e., bins) for any covariate for generating basis functions. If num_knots is a unit-length vector, then the same num_knots are used for each degree (this is not recommended). The default settings for num_knots are recommended, and these defaults decrease num_knots with increasing max_degree and smoothness_orders, which prevents (expensive) combinatorial explosions in the number of higher-degree and higher-order basis functions generated. This allows the complexity of the optimization problem to grow scalably. See details of num_knots more information.

reduce_basis

An optional numeric value bounded in the open unit interval indicating the minimum proportion of 1's in a basis function column needed for the basis function to be included in the procedure to fit the lasso. Any basis functions with a lower proportion of 1's than the cutoff will be removed. Defaults to 1 over the square root of the number of observations. Only applicable for models fit with zero-order splines, i.e. smoothness_orders = 0.

family

A character or a family object (supported by glmnet) specifying the error/link family for a generalized linear model. character options are limited to "gaussian" for fitting a standard penalized linear model, "binomial" for penalized logistic regression, "poisson" for penalized Poisson regression, "cox" for a penalized proportional hazards model, and "mgaussian" for multivariate penalized linear model. Note that passing in family objects leads to slower performance relative to passing in a character family (if supported). For example, one should set family = "binomial" instead of family = binomial() when calling fit_hal.

lambda

User-specified sequence of values of the regularization parameter for the lasso L1 regression. If NULL, the default sequence in cv.glmnet will be used. The cross-validated optimal value of this regularization parameter will be selected with cv.glmnet. If fit_control's cv_select argument is set to FALSE, then the lasso model will be fit via glmnet, and regularized coefficient values for each lambda in the input array will be returned.

id

A vector of ID values that is used to generate cross-validation folds for cv.glmnet. This argument is ignored when fit_control's cv_select argument is FALSE.

weights

observation weights; defaults to 1 per observation.

offset

a vector of offset values, used in fitting.

fit_control

List of arguments, including the following, and any others to be passed to cv.glmnet or glmnet.

- cv_select: A logical specifying if the sequence of specified lambda values should be passed to cv.glmnet in order for a single, optimal value of lambda to be selected according to cross-validation. When cv_select = FALSE, a glmnet model will be used to fit the sequence of (or single) lambda.
- use_min: Specify the choice of lambda to be selected by cv.glmnet. When TRUE, "lambda.min" is used; otherwise, "lambda.1se". Only used when cv_select = TRUE.
• lambda.min.ratio: A glmnet argument specifying the smallest value for lambda, as a fraction of lambda.max, the (data derived) entry value (i.e. the smallest value for which all coefficients are zero). We’ve seen that not setting lambda.min.ratio can lead to no lambda values that fit the data sufficiently well.

• prediction.bounds: An optional vector of size two that provides the lower and upper bounds predictions; not used when family = "cox". When prediction.bounds = "default", the predictions are bounded between min(Y) - sd(Y) and max(Y) + sd(Y) for each outcome (when family = "mgaussian", each outcome can have different bounds). Bounding ensures that there is no extrapolation.

basis_list The full set of basis functions generated from X.
return_lasso A logical indicating whether or not to return the glmnet fit object of the lasso model.
return_x_basis A logical indicating whether or not to return the matrix of (possibly reduced) basis functions used in fit_hal.
yolo A logical indicating whether to print one of a curated selection of quotes from the HAL9000 computer, from the critically acclaimed epic science-fiction film "2001: A Space Odyssey" (1968).

Details

The procedure uses a custom C++ implementation to generate a design matrix of spline basis functions of covariates and interactions of covariates. The lasso regression is fit to this design matrix via cv.glmnet or a custom implementation derived from origami. The maximum dimension of the design matrix is $n \cdot (n \ast 2^d - 1)$, where where $n$ is the number of observations and $d$ is the number of covariates.

For smoothness_orders = 0, only zero-order splines (piece-wise constant) are generated, which assume the true regression function has no smoothness or continuity. When smoothness_orders = 1, first-order splines (piece-wise linear) are generated, which assume continuity of the true regression function. When smoothness_orders = 2, second-order splines (piece-wise quadratic and linear terms) are generated, which assume a the true regression function has a single order of differentiability.

num_knots argument specifies the number of knot points for each covariate and for each max_degree. Fewer knot points can significantly decrease runtime, but might be overly simplistic. When considering smoothness_orders = 0, too few knot points (e.g., < 50) can significantly reduce performance. When smoothness_orders = 1 or higher, then fewer knot points (e.g., 10-30) is actually better for performance. We recommend specifying num_knots with respect to smoothness_orders, and as a vector of length max_degree with values decreasing exponentially. This prevents combinatorial explosions in the number of higher-degree basis functions generated. The default behavior of num_knots follows this logic — for smoothness_orders = 0, num_knots is set to 500/2^j, and for smoothness_orders = 1 or higher, num_knots is set to 200/2^j, where j is the interaction degree. We also include some other suitable settings for num_knots below, all of which are less complex than default num_knots and will thus result in a faster runtime:

• Some good settings for little to no cost in performance:
  - If smoothness_orders = 0 and max_degree = 3, num_knots = c(400, 200, 100).
– If smoothness_orders = 1+ and max_degree = 3, num_knots = c(100, 75, 50).
• Recommended settings for fairly fast runtime:
  – If smoothness_orders = 0 and max_degree = 3, num_knots = c(200, 100, 50).
  – If smoothness_orders = 1+ and max_degree = 3, num_knots = c(50, 25, 15).
• Recommended settings for fast runtime:
  – If smoothness_orders = 0 and max_degree = 3, num_knots = c(100, 50, 25).
  – If smoothness_orders = 1+ and max_degree = 3, num_knots = c(40, 15, 10).
• Recommended settings for very fast runtime:
  – If smoothness_orders = 0 and max_degree = 3, num_knots = c(50, 25, 10).
  – If smoothness_orders = 1+ and max_degree = 3, num_knots = c(25, 10, 5).

Value
Object of class hal9001, containing a list of basis functions, a copy map, coefficients estimated for basis functions, and timing results (for assessing computational efficiency).

Examples
n <- 100
p <- 3
x <- xmat <- matrix(rnorm(n * p), n, p)
y_prob <- plogis(3 * sin(x[, 1]) + sin(x[, 2]))
y <- rbinom(n = n, size = 1, prob = y_prob)
hal_fit <- fit_hal(X = x, Y = y, family = "binomial")
preds <- predict(hal_fit, new_data = x)
HAL Formula term: Generate a single term of the HAL basis

**Description**

HAL Formula term: Generate a single term of the HAL basis

**Usage**

```R
h(
  ..., 
  k = NULL, 
  s = NULL, 
  pf = 1, 
  monotone = c("none", "i", "d"), 
  . = NULL, 
  dot_args_as_string = FALSE, 
  X = NULL 
)
```

**Arguments**

- `...` Variables for which to generate multivariate interaction basis function where the variables can be found in a matrix `X` in a parent environment/frame. Note, just like standard `formula` objects, the variables should not be characters (e.g. do `h(W1,W2)` not `h("W1", "W2")`) `h(W1,W2,W3)` will generate three-way HAL basis functions between `W1`, `W2`, and `W3`. It will not generate the lower dimensional basis functions.

- `k` The number of knots for each univariate basis function used to generate the tensor product basis functions. If a single value then this value is used for the univariate basis functions for each variable. Otherwise, this should be a variable named list that specifies for each variable how many knots points should be used. `h(W1,W2,W3, k = list(W1 = 3, W2 = 2, W3=1))` is equivalent to first binning the variables `W1`, `W2` and `W3` into 3, 2 and 1 unique values and then calling `h(W1,W2,W3)`. This coarsening of the data ensures that fewer basis functions are generated, which can lead to substantial computational speed-ups. If not provided and the variable `num_knots` is in the parent environment, then `s` will be set to `num_knots`.

- `s` The smoothness_orders for the basis functions. The possible values are 0 for piece-wise constant zero-order splines or 1 for piece-wise linear first-order splines. If not provided and the variable `smoothness_orders` is in the parent environment, then `s` will be set to `smoothness_orders`.

- `pf` A penalty.factor value the generated basis functions that is used by `glmnet` in the LASSO penalization procedure. `pf = 1` (default) is the standard penalization factor used by `glmnet` and `pf = 0` means the generated basis functions are unpenalized.
Whether the basis functions should enforce monotonicity of the interaction term. If \( \text{code(s)} = 0 \), this is monotonicity of the function, and, if \( \text{code(s)} = 1 \), this is monotonicity of its derivative (e.g., enforcing a convex fit). Set "none" for no constraints, "i" for a monotone increasing constraint, and "d" for a monotone decreasing constraint. Using "i" constrains the basis functions to have positive coefficients in the fit, and "d" constrains the basis functions to have negative coefficients.

Just like with \texttt{formula} and \texttt{...}, as in \texttt{h(.)} or \texttt{h(...)} is treated as a wildcard variable that generates terms using all variables in the data. The argument \texttt{...} should be a character vector of variable names that \texttt{...} iterates over. Specifically, \texttt{h(..., k=1, . = c("W1", "W2", "W3"))} is equivalent to \texttt{h(W1, k=1) + h(W2, k=1) + h(W3, k=1)}.

\texttt{dot_args_as_string} Whether the arguments \texttt{...} are characters or character vectors and should thus be evaluated directly. When \texttt{TRUE}, the expression \texttt{h("W1", "W2")} can be used.

\texttt{X} An optional design matrix where the variables given in \texttt{...} can be found. Otherwise, \texttt{X} is taken from the parent environment.

---

\textbf{hal9000} \hspace{1cm} \textit{HAL 9000 Quotes}

\textbf{Description}

Prints a quote from the HAL 9000 robot from 2001: A Space Odyssey

\textbf{Usage}

\texttt{hal9000()}

---

\textbf{hal9001} \hspace{1cm} \textit{hal9001}

\textbf{Description}

Package for fitting the Highly Adaptive LASSO (HAL) estimator
**hal_quotes**

**Hal9000 Quotes from "2001: A Space Odyssey"**

**Description**

Curated selection of quotes from the HAL9000 computer, from the critically acclaimed epic science-fiction film "2001: A Space Odyssey" (1968).

**Usage**

hal_quotes

**Format**

A vector of quotes.

**index_first_copy**

Find Copies of Columns

**Description**

Index vector that, for each column in X, indicates the index of the first copy of that column.

**Usage**

index_first_copy(X)

**Arguments**

X Sparse matrix containing columns of indicator functions.

**make_basis_list**

Sort Basis Functions

**Description**

Build a sorted list of unique basis functions based on columns, where each basis function is a list.

**Usage**

make_basis_list(X_sub, cols, order_map)
Arguments

X_sub A subset of the columns of X, the original design matrix.
cols An index of the columns that were reduced to by sub-setting.
order_map A vector with length the original unsubsetted matrix X which specifies the smoothness of the function in each covariate.

Details

Note that sorting of columns is performed such that the basis order equals cols.length() and each basis function is a list(cols, cutoffs).

make_copy_map Build Copy Maps

Description

Build Copy Maps

Usage

make_copy_map(x_basis)

Arguments

x_basis A design matrix consisting of basis (indicator) functions for covariates (X) and terms for interactions thereof.

Value

A list of numeric vectors indicating indices of basis functions that are identical in the training set.

Examples

gedata <- function(n) {
  W1 <- runif(n, -3, 3)
  W2 <- rnorm(n)
  W3 <- runif(n)
  W4 <- rnorm(n)
  g0 <- plogis(0.5 * (-0.8 * W1 + 0.39 * W2 + 0.08 * W3 - 0.12 * W4))
  A <- rbinom(n, 1, g0)
  Q0 <- plogis(0.15 * (2 * A + 2 * A * W1 + 6 * A * W3 * W4 - 3))
  Y <- rbinom(n, 1, Q0)
  data.frame(A, W1, W2, W3, W4, Y)
}
set.seed(1234)
data <- gedata(100)
covars <- setdiff(names(data), "Y")
X <- as.matrix(data[, covars, drop = FALSE])
basis_list <- enumerate_basis(X)
x_basis <- make_design_matrix(X, basis_list)
copy_map <- make_copy_map(x_basis)

make_design_matrix  
Build HAL Design Matrix

Description

Make a HAL design matrix based on original design matrix X and a list of basis functions in argument blist

Usage

make_design_matrix(X, blist, p_reserve = 0.5)

Arguments

X  
Matrix of covariates containing observed data in the columns.

blist  
List of basis functions with which to build HAL design matrix.

p_reserve  
Sparse matrix pre-allocation proportion. Default value is 0.5. If one expects a dense HAL design matrix, it is useful to set p_reserve to a higher value.

Value

A dgCMatrix sparse matrix of indicator basis functions corresponding to the design matrix in a zero-order highly adaptive lasso.

Examples

gedata <- function(n) {
  W1 <- runif(n, -3, 3)
  W2 <- rnorm(n)
  W3 <- runif(n)
  W4 <- rnorm(n)
  g0 <- plogis(0.5 * (-0.8 * W1 + 0.39 * W2 + 0.08 * W3 - 0.12 * W4))
  A <- rbinom(n, 1, g0)
  Q0 <- plogis(0.15 * (2 * A + 2 * A * W1 + 6 * A * W3 * W4 - 3))
  Y <- rbinom(n, 1, Q0)
  data.frame(A, W1, W2, W3, W4, Y)
}
set.seed(1234)
data <- gedata(100)
covars <- setdiff(names(data), "Y")
X <- as.matrix(data[, covars, drop = FALSE])
basis_list <- enumerate_basis(X)
x_basis <- make_design_matrix(X, basis_list)

make_reduced_basis_map

 massa-based reduction of basis functions 

Description

A helper function that finds which basis functions to keep (and equivalently which to discard) based on the proportion of 1's (observations, i.e., "mass") included in a given basis function.

Usage

make_reduced_basis_map(x_basis, reduce_basis_crit)

Arguments

x_basis
A matrix of basis functions with all redundant basis functions already removed.

reduce_basis_crit
A scalar numeric value bounded in the open interval (0,1) indicating the minimum proportion of 1's in a basis function column needed for the basis function to be included in the procedure to fit the Lasso. Any basis functions with a lower proportion of 1's than the specified cutoff will be removed. This argument defaults to NULL, in which case all basis functions are used in the lasso-fitting stage of the HAL algorithm.

Value

A binary numeric vector indicating which columns of the matrix of basis functions to keep (given a one) and which to discard (given a zero).

meets_basis

Compute Values of Basis Functions

Description

Computes and returns the indicator value for the basis described by cols and cutoffs for a given row of X.

Usage

meets_basis(X, row_num, cols, cutoffs, orders)
**predict.hal9001**

**Arguments**

- **X**
  The design matrix, containing the original data.

- **row_num**
  Number for a row index over which to evaluate.

- **cols**
  Numeric for the column indices of the basis function.

- **cutoffs**
  Numeric providing thresholds.

- **orders**
  Numeric providing smoothness orders

---

**predict.hal9001**  
*Prediction from HAL fits*

---

**Description**

Prediction from HAL fits

**Usage**

```r
## S3 method for class 'hal9001'
predict(  
  object,  
  new_data,  
  new_X_unpenalized = NULL,  
  offset = NULL,  
  type = c("response", "link"),  
  ...  
)
```

**Arguments**

- **object**
  An object of class hal9001, containing the results of fitting the Highly Adaptive Lasso, as produced by `fit_hal`.

- **new_data**
  A matrix or data.frame containing new data (i.e., observations not used for fitting the hal9001 object that's passed in via the object argument) for which the hal9001 object will compute predicted values.

- **new_X_unpenalized**
  If the user supplied X_unpenalized during training, then user should also supply this matrix with the same number of observations as new_data.

- **offset**
  A vector of offsets. Must be provided if provided at training.

- **type**
  Either "response" for predictions of the response, or "link" for un-transformed predictions (on the scale of the link function).

- **...**
  Additional arguments passed to predict as necessary.

**Details**

Method for computing and extracting predictions from fits of the Highly Adaptive Lasso estimator, returned as a single S3 objects of class hal9001.
Value

A numeric vector of predictions from a `hal9001` object.

Note

This prediction method does not function similarly to the equivalent method from `glmnet`. In particular, this procedure will not return a subset of lambdas originally specified in calling `fit_hal` nor result in re-fitting. Instead, it will return predictions for all of the lambdas specified in the call to `fit_hal` that constructs object, when `fit_control`'s `cv_select` is set to `FALSE`. When `fit_control`'s `cv_select` is set to `TRUE`, predictions will only be returned for the value of lambda selected by cross-validation.

Description

Predict method for objects of class `SL.hal9001`

Usage

```R
# S3 method for class 'SL.hal9001'
predict(object, newdata, ...)
```

Arguments

- `object` A fitted object of class `hal9001`.
- `newdata` A matrix of new observations on which to obtain predictions.
- `...` Not used.

Value

A numeric vector of predictions from a `SL.hal9001` object based on the provide `newdata`. 
print.formula_hal9001  

Print formula_hal9001 object

Description

Print formula_hal9001 object

Usage

```r
## S3 method for class 'formula_hal9001'
print(x, ...)
```

Arguments

- `x`: A `formula_hal9001` object.
- `...`: Other arguments (ignored).

print.summary.hal9001  

Print Method for Summary Class of HAL fits

Description

Print Method for Summary Class of HAL fits

Usage

```r
## S3 method for class 'summary.hal9001'
print(x, length = NULL, ...)
```

Arguments

- `x`: An object of class `summary.hal9001`.
- `length`: The number of ranked coefficients to be summarized.
- `...`: Other arguments (ignored).
Description

Wrapper for SuperLearner for objects of class hal9001

Usage

\[
\text{SL.hal9001}(Y, X, newX, family, obsWeights, id, max_degree = 2, smoothness_orders = 1, num_knots = 5, \ldots)
\]

Arguments

- **Y**: A numeric vector of observations of the outcome variable.
- **X**: An input matrix with dimensions number of observations -by- number of covariates that will be used to derive the design matrix of basis functions.
- **newX**: A matrix of new observations on which to obtain predictions. The default of NULL computes predictions on training inputs \(X\).
- **family**: A family object (one that is supported by \texttt{glmnet}) specifying the error/link family for a generalized linear model.
- **obsWeights**: A numeric vector of observational-level weights.
- **id**: A numeric vector of IDs.
- **max_degree**: The highest order of interaction terms for which basis functions ought to be generated.
- **smoothness_orders**: An integer vector of length 1 or greater, specifying the smoothness of the basis functions. See the argument smoothness_orders of \texttt{fit_hal} for more information.
- **num_knots**: An integer vector of length 1 or max_degree, specifying the maximum number of knot points (i.e., bins) for each covariate for generating basis functions. See num_knots argument in \texttt{fit_hal} for more information.
- **...**: Additional arguments to \texttt{fit_hal}.
Value

An object of class SL.hal9001 with a fitted hal9001 object and corresponding predictions based on the input data.

Description

Reduce footprint by dropping basis functions with coefficients of zero

Usage

squash_hal_fit(object)

Arguments

object

An object of class hal9001, containing the results of fitting the Highly Adaptive LASSO, as produced by a call to fit_hal.

Value

Object of class hal9001, similar to the input object but reduced such that coefficients belonging to bases with coefficients equal to zero removed.

Examples

# generate simple test data
n <- 100
p <- 3
x <- matrix(rnorm(n * p), n, p)
y <- sin(x[, 1]) * sin(x[, 2]) + rnorm(n, mean = 0, sd = 0.2)

# fit HAL model and squash resulting object to reduce footprint
hal_fit <- fit_hal(X = x, Y = y, yolo = FALSE)
squashed <- squash_hal_fit(hal_fit)
Summary Method for HAL fit objects

Usage

```r
## S3 method for class 'hal9001'
summary(
  object,
  lambda = NULL,
  only_nonzero_coefs = TRUE,
  include_redundant_terms = FALSE,
  round_cutoffs = 3,
  ...
)
```

Arguments

- **object**: An object of class `hal9001`, containing the results of fitting the Highly Adaptive Lasso, as produced by `fit_hal`.
- **lambda**: Optional numeric value of the lambda tuning parameter, for which corresponding coefficient values will be summarized. Defaults to `fit_hal`'s optimal value, `lambda_star`, or the minimum value of `lambda_star`.
- **only_nonzero_coefs**: A logical specifying whether the summary should include only terms with non-zero coefficients.
- **include_redundant_terms**: A logical specifying whether the summary should remove so-called "redundant terms". We define a redundant term (say x1) as a term (1) with basis function corresponding to an existing basis function, a duplicate; and (2) the duplicate contains the x1 term as part of its term, so that x1 terms inclusion would be redundant. For example, say the same coefficient corresponds to these three terms: (1) "I(age >= 50)*I(bmi >= 18)", (2) "I(age >= 50)", and (3) "I(education >= 16)". When `include_redundant_terms` is `FALSE` (default), the second basis function is omitted.
- **round_cutoffs**: An integer indicating the number of decimal places to be used for rounding cutoff values in the term. For example, if "bmi" was numeric that was rounded to the third decimal, in the example above we would have needed to specify `round_cutoffs = 0` in order to yield a term like "I(bmi >= 18)" opposed to something like "I(bmi >= 18.111)". This rounding is intended to simplify the term-wise part of the output and only rounds the basis cutoffs, the `hal9001` model's coefficients are not rounded.
- **...**: Additional arguments passed to `summary`, not supported.
Details

Method for summarizing the coefficients of the Highly Adaptive Lasso estimator in terms of the basis functions corresponding to covariates and interactions of covariates, returned as a single S3 object of class `hal9001`.

Due to the nature of the basis function terms, the summary tables can be extremely wide. The R environment might not be the optimal location to view the summary. Tables can be exported from R to LaTeX with `xtable` package (or similar). Here's an example: `print(xtable(summary(fit)$table, type = "latex"), file = "dt.tex").`

Value

A list summarizing a `hal9001` object's coefficients.
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