

# Package: hierNest (via r-universe)

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**Type** Package

**Title** Penalized Regression with Hierarchical Nested Parameterization Structure

**Version** 1.0.2

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**Description** Efficient implementation of penalized regression with hierarchical nested parametrization for grouped data. The package provides penalized regression methods that decompose subgroup specific effects into shared global effects, Major subgroup specific effects, and Minor subgroup specific effects, enabling structured borrowing of information across related clinical subgroups. Both lasso and hierarchical overlapping group lasso penalties are supported to encourage sparsity while respecting the nested subgroup structure. Efficient computation is achieved through a modified design matrix representation and a custom algorithm for overlapping group penalties.

**License** GPL ( $\geq 2$ )

**Depends** R ( $\geq 3.5$ ),

**Imports** cli, dotCall64, ggplot2, magrittr, Matrix, methods, rlang ( $\geq 1.1.4$ ), RSpectra, tidyr ( $\geq 1.3.0$ ), rTensor, pROC, plotly

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**BugReports** <https://github.com/ZirenJiang/hierNest/issues>

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coef.cv.hierNest

*Extract subgroup-specific coefficients from a cv.hierNest object*

---

## Description

Computes the composite subgroup-specific coefficient vectors from a cross-validated hierarchical nested model. Each subgroup's coefficient is the sum of the overall mean effect, its group (MDC) effect, and its subgroup (DRG) effect:  $\beta^{(s)} = \beta^{overall} + \beta^{group} + \beta^{subgroup}$ .

**Usage**

```
## S3 method for class 'cv.hierNest'
coef(
  object,
  s = c("lambda.min", "lambda.1se"),
  type = c("subgroup", "hierarchical"),
  ...
)
```

**Arguments**

|        |   |
|--------|---|
| object | A fitted <code>cv.hierNest</code> object.   |
| s      | Value of the penalty parameter $\lambda$ at which coefficients are extracted. Either "lambda.min" (default, the $\lambda$ that minimizes CV error) or "lambda.1se" (largest $\lambda$ within 1 SE of the minimum).  |
| type   | Character string specifying the format of the returned coefficients. "subgroup" (default) returns the composite subgroup-specific coefficients (one row per subgroup). "hierarchical" returns the raw hierarchical parameterization (overall, group, and subgroup effect rows). |
| ...    | Not used.   |

**Value**

A numeric matrix of coefficients. Columns correspond to covariates (including "Intercept"). For type = "subgroup", rows are named by subgroup ID. For type = "hierarchical", rows are labeled with overall, group, and subgroup identifiers.

**See Also**

[cv.hierNest](#), [plot\\_contribution](#)

---

|                  |  |
|------------------|--|
| coef.cv.sparsegl | <i>Extract coefficients from a 'cv.sparsegl' object.</i> |
|------------------|--|

---

**Description**

This function extracts coefficients from a cross-validated [`sparsegl()`] model, using the stored "`sparsegl.fit`" object, and the optimal value chosen for '`lambda`'.

**Usage**

```
## S3 method for class 'cv.sparsegl'
coef(object, s = c("lambda.1se", "lambda.min"), ...)
```

**Arguments**

|        |   |
|--------|---|
| object | Fitted [cv.sparsegl()] object.  |
| s      | Value(s) of the penalty parameter ‘lambda’ at which coefficients are desired. Default is the single value ‘s = "lambda.1se"’ stored in the CV object (corresponding to the largest value of ‘lambda’ such that CV error estimate is within 1 standard error of the minimum). Alternatively ‘s = "lambda.min"’ can be used (corresponding to the minimum of cross validation error estimate). If ‘s’ is numeric, it is taken as the value(s) of ‘lambda’ to be used. |
| ...    | Not used.   |

**Value**

The coefficients at the requested value(s) for ‘lambda’.

**See Also**

[cv.sparsegl()] and [predict.cv.sparsegl()].

---

|               |   |
|---------------|---|
| coef.sparsegl | <i>Extract model coefficients from a ‘sparsegl’ object.</i> |
|---------------|---|

---

**Description**

Computes the coefficients at the requested value(s) for ‘lambda’ from a [sparsegl()] object.

**Usage**

```
## S3 method for class 'sparsegl'
coef(object, s = NULL, ...)
```

**Arguments**

|        |  |
|--------|--|
| object | Fitted [sparsegl()] object.  |
| s      | Value(s) of the penalty parameter ‘lambda’ at which coefficients are required. Default is the entire sequence. |
| ...    | Not used.  |

**Details**

‘s’ is the new vector of ‘lambda’ values at which predictions are requested. If ‘s’ is not in the lambda sequence used for fitting the model, the ‘coef’ function will use linear interpolation to make predictions. The new values are interpolated using a fraction of coefficients from both left and right ‘lambda’ indices.

**Value**

The coefficients at the requested values for ‘lambda’.

**See Also**

[sparsegl()] and [predict.sparsegl()].

---

|             |   |
|-------------|---|
| cv.hierNest | <i>Cross-validated hierarchical nested regularization for subgroup models</i> |
|-------------|---|

---

**Description**

Fits regularization paths for hierarchical subgroup-specific penalized learning problems, leveraging nested group structure (such as Major Diagnostic Categories [MDC] and Diagnosis-Related Groups [DRG]) with options for lasso or overlapping group lasso penalties. Performs cross-validation to select tuning parameters for penalization and subgroup structure.

This function enables information sharing across related subgroups by reparameterizing covariate effects into overall, group-specific, and subgroup-specific components and supports structured shrinkage through hierarchical regularization. Users may select between the lasso (hierNest-Lasso) and overlapping group lasso (hierNest-OGLasso) frameworks, as described in Jiang et al. (2024, submitted, see details below).

**Usage**

```
cv.hierNest(
  x,
  y,
  group = NULL,
  family = c("gaussian", "binomial"),
  nlambdas = 100,
  lambda.factor = NULL,
  pred.loss = c("default", "mse", "deviance", "mae", "misclass", "ROC"),
  lambda = NULL,
  pf_group = NULL,
  pf_sparse = NULL,
  intercept = FALSE,
  asparse1 = c(0.5, 20),
  asparse2 = c(0.01, 0.2),
  asparse1_num = 4,
  asparse2_num = 4,
  standardize = TRUE,
  lower_bnd = -Inf,
  upper_bnd = Inf,
  eps = 1e-08,
  maxit = 3e+06,
  hier_info = NULL,
  method = "overlapping",
  partition = "subgroup",
  cvmethod = "general"
)
```

**Arguments**

|               |   |
|---------------|---|
| x             | Matrix of predictors, of dimension $n \times p$ ; each row is an observation. Can be a dense or sparse matrix.  |
| y             | Response variable. For 'family="gaussian"', should be numeric. For 'family="binomial"', should be a factor with two levels or a numeric vector with two unique values.        |
| group         | Optional vector or factor indicating group assignments for variables. Used for custom grouping.   |
| family        | Character string specifying the model family. Options are "gaussian" (default) for least-squares regression, or "binomial" for logistic regression.                           |
| nlambda       | Number of lambda values to use for regularization path. Default is 100.   |
| lambda.factor | Factor determining the minimal value of lambda in the sequence, where 'min(lambda) = lambda.factor * max(lambda)'. See Details.   |
| pred.loss     | Character string indicating loss to minimize during cross-validation. Options include "default", "mse", "deviance", "mae", "misclass", and "ROC".                             |
| lambda        | Optional user-supplied sequence of lambda values (overrides 'nlambda'/'lambda.factor').   |
| pf_group      | Optional penalty factors on the groups, as a numeric vector. Default adjusts for group size.  |
| pf_sparse     | Optional penalty factors on the l1-norm (for sparsity), as a numeric vector.  |
| intercept     | Logical; whether to include an intercept in the model. Default is TRUE.   |
| asparsel      | Relative weight(s) for the first (e.g., group) layer of the overlapping group lasso penalty. Default is c(0.5, 20).   |
| asparsel2     | Relative weight(s) for the second (e.g., subgroup) layer. Default is c(0.01, 0.2).  |
| asparsel1_num | Number of values in asparsel1 grid (for grid search). Default is 4.   |
| asparsel2_num | Number of values in asparsel2 grid (for grid search). Default is 4.   |
| standardize   | Logical; whether to standardize predictors prior to model fitting. Default is TRUE.   |
| lower_bnd     | Lower bound(s) for coefficient values. Default is -Inf.   |
| upper_bnd     | Upper bound(s) for coefficient values. Default is Inf.  |
| eps           | Convergence tolerance for optimization. Default is 1e-8.  |
| maxit         | Maximum number of optimization iterations. Default is 3e6.  |
| hier_info     | Required for 'method = "overlapping"'; a matrix describing the hierarchical structure of the subgroups (see Details).   |
| method        | Character; either "overlapping" for overlapping group lasso, "sparsegl" for sparse group lasso, or "general" for other hierarchical regularization. Default is "overlapping". |
| partition     | Character string; determines subgroup partitioning. Default is "subgroup".  |
| cvmethod      | Cross-validation method. Options include "general" (default), "grid_search", or "user_supply" (for user-supplied grid).   |

## Details

The hierarchical nested framework decomposes covariate effects into overall, group, and subgroup-specific components, with regularization encouraging fusion or sparsity across these hierarchical levels. The function can fit both the lasso penalty (allowing arbitrary zero/non-zero patterns) and the overlapping group lasso penalty (enforcing hierarchical selection structure), as described in Jiang et al. (2024, submitted).

The argument 'hier\_info' must be supplied for "overlapping" method, and encodes the hierarchical relationship between groups and subgroups (e.g., MDCs and DRGs).

Cross-validation is used to select tuning parameters, optionally over a grid for hierarchical penalty weights (asparsel, asparsel2), and the regularization parameter lambda.

## Value

An object containing the fitted hierarchical model and cross-validation results, including:

|             |  |
|-------------|--|
| fit         | Fitted model object.   |
| lambda      | Sequence of lambda values considered.                                  |
| cv_error    | Cross-validation error/loss for each combination of tuning parameters. |
| best_params | Best tuning parameters selected.                                       |
| ...         | Additional diagnostic and output fields.                               |

## References

Jiang, Z., Huo, L., Hou, J., & Huling, J. D. "Heterogeneous readmission prediction with hierarchical effect decomposition and regularization".

---

|               |  |
|---------------|--|
| estimate_risk | <i>Calculate information criteria.</i> |
|---------------|--|

---

## Description

This function uses the degrees of freedom to calculate various information criteria. This function uses the "unknown variance" version of the likelihood. Only implemented for Gaussian regression. The constant is ignored (as in [stats::extractAIC()]).

## Usage

```
estimate_risk(object, x, type = c("AIC", "BIC", "GCV"), approx_df = FALSE)
```

**Arguments**

|           |  |
|-----------|--|
| object    | fitted object from a call to <code>[sparsegl()]</code> .   |
| x         | Matrix. The matrix of predictors used to estimate the ‘sparsegl’ object. May be missing if ‘approx_df = TRUE’.   |
| type      | one or more of AIC, BIC, or GCV.   |
| approx_df | the ‘df’ component of a ‘sparsegl’ object is an approximation (albeit a fairly accurate one) to the actual degrees-of-freedom. However, the exact value requires inverting a portion of ‘X’X’. So this computation may take some time (the default computes the exact df). |

**Value**

a ‘data.frame’ with as many rows as ‘object\$lambda’. It contains columns ‘lambda’, ‘df’, and the requested risk types.

**References**

Vaiter S, Deledalle C, Peyré G, Fadili J, Dossal C. (2012). *The Degrees of Freedom of the Group Lasso for a General Design*. <https://arxiv.org/pdf/1212.6478.pdf>.

**See Also**

`[sparsegl()]` method.

---

example\_data

*Example dataset*

---

**Description**

Example dataset for the sample code of hierNest

**Usage**

```
data(example_data)
```

**Format**

A data frame with X rows and Y columns:

**X** Matrix of predictors

**Y** Outcome variable

**hier\_info** Group information passed to hierNest functions

---

 hierNest

*Fit Hierarchical Nested Regularization Model (hierNest)*


---

### Description

Fits a hierarchical nested penalized regression model for subgroup-specific effects using overlapping group lasso penalties. This function encodes the hierarchical structure (e.g., MDC and DRG) via a reparameterized design matrix and enables information borrowing across related subgroups.

### Usage

```

hierNest(
  x,
  y,
  group = NULL,
  family = c("gaussian", "binomial"),
  nlambda = 100,
  lambda.factor = NULL,
  lambda = NULL,
  pf_group = NULL,
  pf_sparse = NULL,
  intercept = FALSE,
  asparse1 = 1,
  asparse2 = 0.05,
  standardize = TRUE,
  lower_bnd = -Inf,
  upper_bnd = Inf,
  eps = 1e-08,
  maxit = 3e+06,
  hier_info = NULL,
  random_asparse = FALSE,
  method = "overlapping"
)

```

### Arguments

|               |  |
|---------------|--|
| x             | Matrix of predictors ( $n \times p$ ), where each row is an observation.                 |
| y             | Response variable (numeric for "gaussian", binary or factor for "binomial").             |
| group         | Optional grouping vector (not required for "overlapping" method).                        |
| family        | Model family; either "gaussian" for least squares or "binomial" for logistic regression. |
| nlambda       | Number of lambda values in the regularization path (default: 100).                       |
| lambda.factor | Factor for minimal value of lambda in the sequence.                                      |
| lambda        | Optional user-supplied lambda sequence.  |
| pf_group      | Penalty factor(s) for each group; defaults to sqrt(group size).                          |

|                 |  |
|-----------------|--|
| pf_sparse       | Penalty factors for individual predictors (L1 penalty).                                  |
| intercept       | Logical; should an intercept be included? Default is FALSE.                              |
| asparsel        | Relative weight for group-level penalty (default: 1).                                    |
| asparsel2       | Relative weight for subgroup-level penalty (default: 0.05).                              |
| standardize     | Logical; standardize predictors? Default is TRUE.  |
| lower_bnd       | Lower bound for coefficients (default: -Inf).  |
| upper_bnd       | Upper bound for coefficients (default: Inf).   |
| eps             | Convergence tolerance (default: 1e-8).   |
| maxit           | Maximum number of optimization iterations (default: 3e6).                                |
| hier_info       | Required. Matrix encoding the hierarchical structure (see Details).                      |
| random_asparsel | Logical; use random sparse penalty? Default: FALSE.                                      |
| method          | Type of hierarchical regularization ("overlapping" [default], "sparsegl", or "general"). |

### Details

This function builds a hierarchical design matrix reflecting group/subgroup structure (e.g., Major Diagnostic Categories [MDCs] and Diagnosis Related Groups [DRGs]), encoding overall, group-specific, and subgroup-specific effects. It fits a penalized model using overlapping group lasso, as described in Jiang et al. (2024, submitted). The main computational engine is `hierNest::overlapping_gl`.

### Value

Returns a model fit object as produced by `hierNest::overlapping_gl`, including selected coefficients, cross-validation results, and tuning parameters.

### References

Jiang, Z., Huo, L., Hou, J., Vaughan-Sarrazin, M., Smith, M. A., & Huling, J. D. (2024). Heterogeneous readmission prediction with hierarchical effect decomposition and regularization.

---

make\_irls\_warmup

*Create starting values for iterative reweighted least squares*

---

### Description

This function may be used to create potentially valid starting values for calling `[sparsegl()]` with a `[stats::family()]` object. It is not typically necessary to call this function (as it is used internally to create some), but in some cases, especially with custom generalized linear models, it may improve performance.

### Usage

```
make_irls_warmup(nobs, nvars, b0 = 0, beta = double(nvars), r = double(nobs))
```

**Arguments**

|       |   |
|-------|---|
| nobs  | Number of observations in the response (or rows in 'x').                                |
| nvars | Number of columns in 'x'  |
| b0    | Scalar. Initial value for the intercept.  |
| beta  | Vector. Initial values for the coefficients. Must be length 'nvars' (or a scalar).      |
| r     | Vector. Initial values for the deviance residuals. Must be length 'nobs' (or a scalar). |

**Details**

Occasionally, the irls fitting routine may fail with an admonition to create valid starting values.

**Value**

List of class 'irlsspgl\_warmup'

---

|                |   |
|----------------|---|
| overlapping_gl | <i>Fit an Overlapping Group Lasso Model</i> |
|----------------|---|

---

**Description**

This function fits an overlapping group lasso model with hierarchical regularization, allowing both sparsity within groups and across groups.

**Usage**

```
overlapping_gl(
  x,
  y,
  group = NULL,
  family = c("gaussian", "binomial"),
  nlambdas = 100,
  lambda.factor = ifelse(nobs < nvars, 0.01, 1e-04),
  lambda = NULL,
  pf_group = sqrt(bs),
  pf_sparse = rep(1, nvars),
  intercept = TRUE,
  asparse1 = 1,
  asparse2 = 0.05,
  standardize = TRUE,
  lower_bnd = -Inf,
  upper_bnd = Inf,
  weights = NULL,
  offset = NULL,
  warm = NULL,
  trace_it = 0,
```

```

dfmax = as.integer(max(group)) + 1L,
pmax = min(dfmax * 1.2, as.integer(max(group))),
eps = 1e-08,
maxit = 3e+06,
cn,
drnix,
drnix,
cn_s,
cn_e,
random_asparse = FALSE
)

```

### Arguments

|                            |   |
|----------------------------|---|
| <code>x</code>             | A numeric matrix of predictor variables (no missing values allowed).  |
| <code>y</code>             | A numeric vector of response variable values.   |
| <code>group</code>         | An integer vector defining the group membership for each predictor. Default is NULL (each predictor forms its own group). |
| <code>family</code>        | A character string specifying the model family. Options are "gaussian" (default) and "binomial".                          |
| <code>nlambda</code>       | Number of lambda values. Default is 100.  |
| <code>lambda.factor</code> | Factor determining minimum lambda as a fraction of maximum lambda. Default depends on dimensionality.                     |
| <code>lambda</code>        | Numeric vector of lambda values. If provided, overrides 'nlambda' and 'lambda.factor'.                                    |
| <code>pf_group</code>      | Penalty factor for groups. Default is square root of group size.  |
| <code>pf_sparse</code>     | Penalty factor for individual predictors. Default is 1 for each predictor.  |
| <code>intercept</code>     | Logical; whether to include an intercept. Default is TRUE.  |
| <code>asparsel</code>      | Sparsity penalty factor controlling group-level sparsity. Default is 1.   |
| <code>asparsel2</code>     | Sparsity penalty factor controlling within-group sparsity. Default is 0.05.   |
| <code>standardize</code>   | Logical; if TRUE, standardizes predictors before fitting. Default is TRUE.  |
| <code>lower_bnd</code>     | Numeric vector specifying lower bounds for coefficients. Default is -Inf.   |
| <code>upper_bnd</code>     | Numeric vector specifying upper bounds for coefficients. Default is Inf.  |
| <code>weights</code>       | Optional numeric vector of observation weights. Currently limited functionality.  |
| <code>offset</code>        | Optional numeric vector specifying a known component to be included in the linear predictor.                              |
| <code>warm</code>          | Optional initial values for optimization.   |
| <code>trace_it</code>      | Integer indicating the verbosity level. Default is 0 (no output).   |
| <code>dfmax</code>         | Maximum number of groups allowed in the model. Default derived from groups.   |
| <code>pmax</code>          | Maximum number of predictors allowed in the model. Default derived from groups.   |
| <code>eps</code>           | Numeric convergence threshold for optimization. Default is 1e-08.   |
| <code>maxit</code>         | Maximum number of iterations for optimization. Default is 3e+06.  |

|                |   |
|----------------|---|
| cn             | Additional internal numeric parameter for optimization.                         |
| drgix, drgiy   | Numeric vectors specifying indices for specific group and predictor structures. |
| cn_s, cn_e     | Numeric vectors specifying starting and ending indices for substructures.       |
| random_asparse | Logical; if TRUE, randomly selects sparsity parameters. Default is FALSE.       |

**Value**

An object of class 'sparsegl' containing:

|                     |                                     |
|---------------------|-------------------------------------|
| call                | The matched function call.          |
| lambda              | The lambda values used for fitting. |
| aspase1, aspase2    | Sparsity parameters used.           |
| nobs                | Number of observations.             |
| pf_group, pf_sparse | Penalty factors used.               |
| coefficients        | Estimated coefficients.             |

Additional components relevant to model diagnostics and fitting.

---

|                  |  |
|------------------|--|
| plot.cv.hierNest | <i>Plot method for cv.hierNest objects</i> |
|------------------|--|

---

**Description**

Plot method for cv.hierNest objects

**Usage**

```
## S3 method for class 'cv.hierNest'
plot(x, type = c("coefficients", "Subgroup effects"), ...)
```

**Arguments**

|      |  |
|------|--|
| x    | An object of class cv.hierNest   |
| type | Plot type, e.g. "fit" for observed vs fitted (default), others you may add later (e.g., "coef"). |
| ...  | Other parameters   |

**Value**

Invisibly returns x

---

|                   |  |
|-------------------|--|
| plot_contribution | <i>Boxplot or bar chart of per-covariate contributions to the linear predictor</i> |
|-------------------|--|

---

### Description

For each subject, computes the contribution of each covariate to the linear predictor as  $x_j \cdot \beta_j$ , where  $\beta_j$  is the subject's subgroup-specific coefficient for covariate  $j$ .

When no `subject_id` is supplied, produces a boxplot showing the distribution of contributions across all subjects. When a single `subject_id` is supplied, produces a bar chart for that subject.

### Usage

```
plot_contribution(
  object,
  newx,
  hier_info,
  s = c("lambda.min", "lambda.1se"),
  subject_id = NULL,
  top_n = NULL,
  include_intercept = FALSE
)
```

### Arguments

|                                |  |
|--------------------------------|--|
| <code>object</code>            | A fitted <code>cv.hierNest</code> object.  |
| <code>newx</code>              | Numeric matrix of predictor values ( $n \times p$ ), with the same columns as the original training data.  |
| <code>hier_info</code>         | Numeric matrix encoding the hierarchical structure for the subjects in <code>newx</code> . First column is group (MDC) membership; second column is subgroup (DRG) membership. Must have <code>nrow(newx)</code> rows. |
| <code>s</code>                 | Value of <code>lambda</code> to use for coefficient extraction. Either <code>"lambda.min"</code> (default) or <code>"lambda.1se"</code> .  |
| <code>subject_id</code>        | Optional. A row index (integer) or row name (character) identifying a single subject in <code>newx</code> . If provided, a bar chart of that subject's covariate contributions is returned instead of a boxplot.       |
| <code>top_n</code>             | Optional integer. If provided, only the top <code>top_n</code> covariates (by median absolute contribution across subjects, or by absolute contribution for a single subject) are shown.                               |
| <code>include_intercept</code> | Logical; if TRUE, include the intercept contribution as a separate covariate. Default is FALSE.  |

**Details**

The function uses `coef.cv.hierNest` to obtain the composite subgroup-specific coefficients. Each subject is mapped to their subgroup via `hier_info`, and the element-wise product  $x_j\beta_j$  quantifies how much each covariate contributes to that subject's predicted risk (linear predictor scale).

Covariates are ordered on the x-axis by descending median absolute contribution (boxplot) or descending absolute contribution (bar chart), so the most influential variables appear on the left.

**Value**

A ggplot2 object.

**See Also**

`coef.cv.hierNest`, `cv.hierNest`

---

`predict.cv.sparsegl`    *Make predictions from a 'cv.sparsegl' object.*

---

**Description**

This function makes predictions from a cross-validated `[cv.sparsegl()]` object, using the stored 'sparsegl.fit' object, and the value chosen for 'lambda'.

**Usage**

```
## S3 method for class 'cv.sparsegl'
predict(
  object,
  newx,
  s = c("lambda.1se", "lambda.min"),
  type = c("link", "response", "coefficients", "nonzero", "class"),
  ...
)
```

**Arguments**

|                     |   |
|---------------------|---|
| <code>object</code> | Fitted <code>[cv.sparsegl()]</code> object.   |
| <code>newx</code>   | Matrix of new values for 'x' at which predictions are to be made. Must be a matrix. This argument is mandatory.   |
| <code>s</code>      | Value(s) of the penalty parameter 'lambda' at which coefficients are desired. Default is the single value 's = "lambda.1se"' stored in the CV object (corresponding to the largest value of 'lambda' such that CV error estimate is within 1 standard error of the minimum). Alternatively 's = "lambda.min"' can be used (corresponding to the minimum of cross validation error estimate). If 's' is numeric, it is taken as the value(s) of 'lambda' to be used. |

|      |  |
|------|--|
| type | Type of prediction required. Type "link" gives the linear predictors for "binomial"; for "gaussian" models it gives the fitted values. Type "response" gives predictions on the scale of the response (for example, fitted probabilities for "binomial"); for "gaussian" type "response" is equivalent to type "link". Type "coefficients" computes the coefficients at the requested values for 's'. Type "class" applies only to "binomial" models, and produces the class label corresponding to the maximum probability. Type "nonzero" returns a list of the indices of the nonzero coefficients for each value of s. |
| ...  | Not used.  |

**Value**

A matrix or vector of predicted values.

**See Also**

[cv.sparsegl()] and [coef.cv.sparsegl()].

---

predict.sparsegl      *Make predictions from a 'sparsegl' object.*

---

**Description**

Similar to other predict methods, this function produces fitted values and class labels from a fitted [`'sparsegl'`] object.

**Usage**

```
## S3 method for class 'sparsegl'
predict(
  object,
  newx,
  s = NULL,
  type = c("link", "response", "coefficients", "nonzero", "class"),
  ...
)
```

**Arguments**

|        |  |
|--------|--|
| object | Fitted [sparsegl()] model object.  |
| newx   | Matrix of new values for 'x' at which predictions are to be made. Must be a matrix. This argument is mandatory.                        |
| s      | Value(s) of the penalty parameter 'lambda' at which predictions are required. Default is the entire sequence used to create the model. |

|      |  |
|------|--|
| type | Type of prediction required. Type "link" gives the linear predictors for "binomial"; for "gaussian" models it gives the fitted values. Type "response" gives predictions on the scale of the response (for example, fitted probabilities for "binomial"); for "gaussian" type "response" is equivalent to type "link". Type "coefficients" computes the coefficients at the requested values for 's'. Type "class" applies only to "binomial" models, and produces the class label corresponding to the maximum probability. Type "nonzero" returns a list of the indices of the nonzero coefficients for each value of s. |
| ...  | Not used.  |

### Details

's' is the new vector of 'lambda' values at which predictions are requested. If 's' is not in the lambda sequence used for fitting the model, the 'coef' function will use linear interpolation to make predictions. The new values are interpolated using a fraction of coefficients from both left and right 'lambda' indices.

### Value

The object returned depends on type.

### See Also

[sparsegl()], [coef.sparsegl()].

---

predict\_hierNest      *Predict Method for hierNest Objects*

---

### Description

Provides predictions from a fitted hierarchical model ('hierNest') using new data.

### Usage

```
predict_hierNest(
  object,
  newx,
  hier_info,
  type = c("link", "response", "coefficients", "nonzero", "class"),
  ...
)
```

**Arguments**

|           |  |
|-----------|--|
| object    | A fitted hierNest model object.  |
| newx      | A numeric matrix of new predictor values for prediction.   |
| hier_info | A numeric matrix with hierarchical grouping information. First column is MDC-level grouping; second column is DRG-level grouping.        |
| type      | Character string specifying the type of prediction required. Options include "link", "response", "coefficients", "nonzero", and "class". |
| ...       | Additional arguments passed to lower-level prediction methods.   |

**Details**

This function prepares a hierarchical design matrix based on 'hier\_info', constructs the required Khatri-Rao product, and reorganizes it before generating predictions from the provided 'object'.

**Value**

Predictions based on the specified 'type'. Typically, returns:

- Numeric vector or matrix of predicted values (for "link" or "response").
- Model coefficients (for "coefficients").
- Nonzero coefficient indices (for "nonzero").
- Class labels for categorical outcomes (for "class").

---

zero\_norm

*Calculate common norms*

---

**Description**

Calculate different norms of vectors with or without grouping structures.

**Usage**

zero\_norm(x)

one\_norm(x)

two\_norm(x)

grouped\_zero\_norm(x, gr)

grouped\_one\_norm(x, gr)

grouped\_two\_norm(x, gr)

grouped\_sp\_norm(x, gr, asparse)

```
gr_one_norm(x, gr)
```

```
gr_two_norm(x, gr)
```

```
sp_group_norm(x, gr, aspase = 0.05)
```

### Arguments

|        |   |
|--------|---|
| x      | A numeric vector.   |
| gr     | An integer (or factor) vector of the same length as x.                    |
| aspase | Scalar. The weight to put on the l1 norm when calculating the group norm. |

### Value

A numeric scalar or vector

### Functions

- zero\_norm(): l0-norm (number of nonzero entries).
- one\_norm(): l1-norm (Absolute-value norm).
- two\_norm(): l2-norm (Euclidean norm).
- grouped\_zero\_norm(): A vector of group-wise l0-norms.
- grouped\_one\_norm(): A vector of group-wise l1-norms.
- grouped\_two\_norm(): A vector of group-wise l2-norms.
- grouped\_sp\_norm(): A vector of length 'unique(gr)' consisting of the 'aspase' convex combination of the l1 and l2-norm for each group.
- gr\_one\_norm(): The l1-norm norm of a vector (a scalar).
- gr\_two\_norm(): The sum of the group-wise l2-norms of a vector (a scalar).
- sp\_group\_norm(): The sum of the 'aspase' convex combination of group l1 and l2-norms vectors (a scalar).

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