

Package: adlaplace (via r-universe)

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Type Package

Title Laplace Approximations Using Automatic Differentiation

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Description Computes Laplace approximations for hierarchical models using automatic differentiation (CppAD) and trust-region optimization.

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Contents

.type_factor_levels	3
ad_data	3
ad_data-class	5
ad_fun	5
ad_fun-class	6
ad_fun_ptr	7
ad_fun_ptr-class	8
ad_shards	8
adlaplace	9
adlaplace_cpp	11
adlaplace_fit-methods	12
apply_theta_log	13
binomial-class	14
by_group	15
by_group-class	15
c.ad_fun_ptr	16
clone_ad_fun_ptr	16
collect_terms	17
format_parameters	17
fpoly-class	18
gamm	19
gaussian-class	20
germany	21
has_openmp	22
iid-class	22
inner_opt	23
intercept-class	24
iwp	26
iwp-class	27
laplace_half_H_inv	28
linear-class	28
log_lik_laplace	30
mean_mvquad	31
model-class	32
model-generics	33
model_data	34
nbinom-class	35
outer_optim_wrappers	36
rmvnldl	37
rpoly-class	38

Arguments

y	Response vector (default empty), or an existing ad_data object to recast as another shard.
A	Random-effects design matrix ($\text{nrow}(A) = \text{length}(y)$; any Matrix class).
X	Fixed-effects design matrix ($\text{nrow}(X) = \text{length}(y)$; any Matrix class).
beta_map	Beta parameter map (any Matrix; coerced to ngCMatrix), or a shorthand: length-1 integer n gives <code>Matrix::Matrix(nrow = n, ncol = 0)</code> ; length-2 <code>c(row, n)</code> gives one column with a structural 1 at row row (1-based); and <code>list(indices, n)</code> gives one column per entry in indices with structural 1s at those rows. nrow is the global beta block size; column j maps local X column j to a global beta row. When $\text{ncol}(X) > 0$, $\text{ncol}(\text{beta_map})$ must equal $\text{ncol}(X)$ with exactly one nonzero per column. If NULL, defaults to <code>Matrix::Diagonal(ncol(X))</code> when $\text{ncol}(X) > 0$, otherwise an empty matrix.
gamma_map	Gamma parameter map (any Matrix; coerced to ngCMatrix; at most one nonzero per row and column), or the same shorthand forms as beta_map. Column j maps local A column j into the global gamma block. When $\text{ncol}(A) > 0$, $\text{ncol}(\text{gamma_map})$ must equal $\text{ncol}(A)$ with exactly one nonzero per column. If NULL, defaults to <code>Matrix::Diagonal(ncol(A))</code> when $\text{ncol}(A) > 0$, otherwise an empty matrix.
theta_map	Theta parameter map (any Matrix; coerced to ngCMatrix), or the same shorthand forms as beta_map. Column j selects global theta row for component j. If NULL, defaults to an empty matrix.
elgm_matrix	Optional ELGM map (any Matrix; coerced to ngCMatrix). If NULL, defaults to <code>Matrix::Matrix(nrow = \text{length}(y), ncol = 0)</code> .
ad_fun	Registered AD density name for this shard (optional).
ad_kind	Shard kind ("observations", "parameters", "random"; optional).
package	Package recording AD tapes for this shard (optional).
precision	Optional precision payload (any R object).
weights	Optional per-observation weights (e.g. binomial trial counts). Empty means all ones.
formula	Model formula or named list of term objects.
data	Data frame.
verbose	Print parsing messages.

Value

List with terms, y, A, X, data, info. The `info$theta` data frame includes a logical transform column: TRUE means the parameter is stored on the log scale when `config$transform_theta` is TRUE (default for all rows unless a model term sets otherwise). `info$parameters` stacks beta and theta rows with the same transform column (FALSE for fixed effects); `init`, `lower`, and `upper` are already on the optimization (log) scale where transform is TRUE.

Functions

- `ad_data()`: Construct an ad_data object

ad_data-class	<i>Per-shard layout for AD density evaluation</i>
---------------	---

Description

Per-shard layout for AD density evaluation

Slots

y Response vector.

ATp Transpose of random-effects design matrix.

XTp Transpose of fixed-effects design matrix.

beta_map Beta parameter map (ngCMatrix). nrow is the global beta block size; column j has one structural nonzero at row r meaning local design column j of X uses global beta r. When $\text{ncol}(X) > 0$, require $\text{ncol}(\text{beta_map}) == \text{ncol}(X)$.

gamma_map Gamma parameter map (ngCMatrix); at most one structural nonzero per row and column. Same column convention as beta_map for local A columns into the global gamma block.

theta_map Theta parameter map (ngCMatrix). Column j selects global theta row r for this shard's theta component j.

elgm_matrix Optional exposure-lag map (ngCMatrix; empty by default).

ad_fun Registered AD density name for this shard.

ad_kind Shard kind ("observations", "parameters", "random").

package Package name whose shared library records tapes for this shard (defaults to "adlaplace" when missing).

precision Optional precision payload (any R object).

weights Optional per-observation weights (e.g. binomial trial counts). Empty means all ones.

ad_fun	<i>Build AD function with Hessian templates</i>
--------	---

Description

Build AD function with Hessian templates

Usage

```
ad_fun(x, config = NULL, num_threads = 1L, ...)

## S4 method for signature 'ad_fun_ptr'
ad_fun(x, config = NULL, num_threads = 1L, ...)

## S4 method for signature 'ad_data'
ad_fun(x, config = NULL, num_threads = 1L, ...)

## S4 method for signature 'list'
ad_fun(x, config = NULL, num_threads = 1L, ...)
```

Arguments

x	One of: an ad_fun_ptr, an ad_data shard with ad_kind/ad_fun set, or a model_data() bundle (list with observations, random, parameters).
config	Model configuration list. Required for ad_data and model_data; unused for ad_fun_ptr. For model_data, beta, gamma, and theta are filled from x\$data\$info when omitted (only shards, transform_theta, etc. are needed). When config\$num_threads is set, it overrides the num_threads argument for ad_data and model_data methods.
num_threads	Positive integer; OpenMP thread count for inner_opt and parallel trace_hinv.t. Shards are assigned owner_thread = shard_index % num_threads at attach time. Default 1L (serial).
...	For ad_fun_ptr input, optional additional ad_fun_ptr shards to combine before attaching templates.

Value

Object of class ad_fun.

ad_fun-class	<i>AD function with Hessian templates attached</i>
--------------	--

Description

AD function with Hessian templates attached

Slots

ptr Raw combined handle (ad_fun_ptr).
group_sparsity Per-group inner gradient sparsity patterns.
outer Outer Hessian template (dgCMatrix).
inner Inner-gamma Hessian template (dgCMatrix).
map_outer Outer Hessian shard map.

map_inner Inner Hessian shard map.

parallel_map Thread-affinity map (ngCMatrix); rows are shards, columns are threads, with one nonzero per shard row.

chol_inner Symbolic LDL factor or empty sparse matrix.

chol_inner_list Numeric LDL list for C++ (L1, Linv, perm, perm_inv, half_H_inv, H_inv), plus trace_columns (per-shard column indices for trace_hinv_t).

sizes Named numeric vector beta/gamma/theta.

info List of parameter metadata (beta, gamma, theta, parameters); populated from model_data()\$data\$info when the handle is built from a model-data bundle, otherwise empty.

ad_fun_ptr	<i>Build raw AD handle for one density shard</i>
------------	--

Description

Constructs CppAD tapes for a single shard. The density kind and name come from data@ad_kind and data@ad_fun. For ad_kind = "random", data@precision is passed straight to the backend. For random_diagonal it must be a numeric vector of diagonal precision weights with length == ncol(gamma_map). Missing precision means the shard is not built (e.g. diffuse rpoly with sd = Inf via model_data()); calling ad_fun_ptr() without it is an error.

Usage

```
ad_fun_ptr(data, config)
```

Arguments

data	An ad_data object with ad_kind and ad_fun slots set.
config	Model configuration list (beta, theta, etc.). gamma is optional; when missing, zeros of length nrow(gamma_map) are used as AD tape seeds. config\$num_threads does not assign OpenMP threads; use num_threads on ad_fun after merging shards.

Details

Merge handles for multiple shards with c() before calling [ad_fun](#).

Value

External pointer of class ad_fun_ptr.

ad_fun_ptr-class	<i>Raw AD handle (external pointer)</i>
------------------	---

Description

S4 class for a C++ AD backend handle. Objects are created by `ad_fun_ptr`, combined with `c.ad_fun_ptr`, and typically passed to `ad_fun` to attach Hessian templates.

See Also

`ad_fun_ptr`, `ad_fun`, `clone_ad_fun_ptr`

ad_shards	<i>Partition observations into AD shards by sparsity pattern</i>
-----------	--

Description

The function partitions columns into `num_shards` shards using quantiles of the first right singular vector, optionally using **RSpectra** for efficiency when available.

Arguments

<code>A</code>	Random-effects design matrix (<code>nrow(A)</code> = number of observations).
<code>elgm_matrix</code>	A numeric matrix (or matrix-like object) for extended latent gaussian models.
<code>num_shards</code>	Integer giving the maximum number of shards to construct. The actual number of shards may be smaller if fewer distinct loadings are present.
<code>min_groups</code>	Integer giving the minimum number of shards. When there are fewer distinct singular-vector loadings than <code>min_groups</code> , only the largest exact-loading group is split until <code>min_groups</code> is reached (or that group cannot be split further).

Details

The resulting grouping is returned as a sparse matrix whose columns correspond to shards and whose entries are the singular-vector loadings. Shards are ordered from most heterogeneous to most homogeneous.

If the **RSpectra** package is available, the leading singular vector is computed using `RSpectra::svds`; otherwise, a full singular value decomposition via `svd` is used.

Shard boundaries are defined by empirical quantiles of the loadings. Shards are subsequently re-ordered so that shards with larger within-shard variability appear first.

Value

A sparse matrix of class "dgCMatrix" (from **Matrix**), with one column per shard and one row per observation. Nonzero entries correspond to singular-vector loadings.

Examples

```
set.seed(1)
A <- matrix(rnorm(100), 20, 5)
G <- ad_shards(A, num_shards = 3)
G
```

adlplace

Fit a hierarchical model by Laplace-approximate maximum likelihood

Description

One-call interface to the **adlplace** pipeline: parse the formula with `model_data()`, partition observations into AD shards with `ad_shards()`, build the AD handle with `ad_fun()`, maximize the Laplace-approximate marginal likelihood with `optim()` using the exact profile gradient (`outer_fn` / `outer_gr`), and return a fitted-model object with standard accessor methods.

Usage

```
adlplace(
  formula,
  data,
  config = list(),
  num_threads = 1L,
  num_shards = 100L,
  control = list(),
  control_inner = list(maxit = 100L, report.level = 0, report.freq = 0),
  method = "L-BFGS-B",
  hessian = TRUE,
  verbose = FALSE,
  na_omit = TRUE
)
```

Arguments

formula	Model formula built from model terms, e.g. <code>nbinom(y) ~ x + iid(g)</code> , a list of term objects from <code>collect_terms</code> , or a precomputed <code>model_data</code> result (in which case data is ignored).
data	Data frame containing the variables referenced in formula. Not needed when formula is already a <code>model_data</code> result.
config	List of backend options merged over the defaults <code>list(transform_theta = TRUE, verbose = verbose)</code> . When <code>config\$shards</code> is missing it is filled by <code>ad_shards(A, num_shards = num_shards)</code> .
num_threads	Positive integer, OpenMP thread count for the AD handle.
num_shards	Target number of observation shards for <code>ad_shards()</code> .
control	List of <code>optim</code> control options, merged over the defaults <code>list(maxit = 200L, parscale = <from terms>)</code> .

<code>control_inner</code>	List of control options for the inner (random effects) trust-region optimizer.
<code>method</code>	<code>optim</code> method; bounds are used for "L-BFGS-B" and "Brent" only.
<code>hessian</code>	Logical. When TRUE (default), <code>optim</code> numerically differentiates the exact AD profile gradient at the optimum, giving the observed information used by <code>vcov.adlaplace_fit</code> , <code>summary.adlaplace_fit</code> , and <code>confint.adlaplace_fit</code> .
<code>verbose</code>	Logical, print progress.
<code>na_omit</code>	Passed to <code>model_data()</code> .

Value

An object of class "adlaplace_fit": a list with components

call, formula The matched call and model formula.

par Named outer parameter estimates on the optimization scale (log scale where `par_info$transform` is TRUE).

par_info Parameter metadata table (`info$parameters`).

vcov Inverse of the numerically differentiated Hessian of the negative log likelihood (optimization scale), or NULL.

gamma Named random-effect modes at the optimum.

logLik, nobs Laplace-approximate marginal log likelihood and the number of observations.

optim Full `optim` result, including hessian when requested.

details Result of `log_lik_laplace(deriv = TRUE)` at the optimum (gradient, Cholesky factors, inner solution).

model_data, ad_fun, config, control_inner, cache Objects needed by the accessor methods and for warm restarts.

See Also

`summary.adlaplace_fit`, `predict.adlaplace_fit`, `log_lik_laplace`, `model_data`

Examples

```
## Not run:
fit <- adlaplace(
  nbinom(y, lower = 1e-9, init = 0.15) ~ x + iid(g, init = 0.1),
  data = dat, num_threads = 2L
)
summary(fit)
confint(fit)

## End(Not run)
```

adlaplace_cpp	C++ backend entry points
---------------	--------------------------

Description

Low-level entry points exposed to R. Accept an `ad_fun` S4 object (via `@ptr`) or a raw `ad_fun_ptr`.

Usage

```
joint_log_dens(ad_fun, x, shards = NULL, negative = TRUE)

grad(ad_fun, x, shards = NULL, inner = FALSE, negative = TRUE)

hessian(
  ad_fun,
  x,
  shards = NULL,
  inner = FALSE,
  verbose = FALSE,
  negative = TRUE
)

trace_hinv_t(ad_fun, x, LinvPt, LinvPtColumns, verbose = FALSE)

fun_obj_fdfh(ad_fun, parameters, gamma, inner = TRUE, verbose = FALSE)
```

Arguments

<code>ad_fun</code>	An <code>ad_fun</code> S4 object or <code>ad_fun_ptr</code> handle.
<code>x</code>	Numeric parameter vector of length <code>Nparams</code> .
<code>shards</code>	Optional integer vector of 0-based shard indices; <code>NULL</code> or <code>integer(0)</code> evaluates all shards.
<code>negative</code>	Logical (default <code>TRUE</code>). If <code>TRUE</code> , return the negative log density $-\ell(x)$ and its derivatives (minimization / trustOptim sign, consistent with <code>inner_opt()</code>). If <code>FALSE</code> , return $\ell(x)$, $\nabla\ell$, and $\nabla^2\ell$.
<code>inner</code>	Logical; if <code>TRUE</code> , evaluate inner- γ derivatives; if <code>FALSE</code> , evaluate outer derivatives at the full parameter vector.
<code>verbose</code>	Logical; if <code>TRUE</code> , print threads, shards, and CppAD session phases.
<code>LinvPt, LinvPtColumns</code>	Sparse factors for <code>trace_hinv_t()</code> .
<code>parameters</code>	Numeric vector of length <code>Nbeta + Ntheta</code> .
<code>gamma</code>	Numeric vector of length <code>Ngamma</code> .

Value

See individual functions.

Sign convention

With default `negative = TRUE`, `joint_log_dens()`, `grad()`, and `hessian()` match `inner_opt()` (negative log-density). Set `negative = FALSE` for the joint log density and its derivatives at the same `x`.

adlaplace_fit-methods *Methods for fitted adlaplace models*

Description

Standard accessor methods for objects of class "adlaplace_fit" returned by `adlaplace()`. Estimates are stored on the optimization scale (log scale for parameters with `par_info$transform = TRUE`, e.g. standard deviations); methods with a `transform` argument report the natural scale by exponentiating and, for standard errors, applying the delta method.

Usage

```
## S3 method for class 'adlaplace_fit'
coef(object, transform = TRUE, ...)

## S3 method for class 'adlaplace_fit'
vcov(object, transform = FALSE, ...)

## S3 method for class 'adlaplace_fit'
logLik(object, ...)

## S3 method for class 'adlaplace_fit'
nobs(object, ...)

## S3 method for class 'adlaplace_fit'
confint(object, parm, level = 0.95, transform = TRUE, ...)

## S3 method for class 'adlaplace_fit'
fitted(object, ...)

## S3 method for class 'adlaplace_fit'
predict(object, newdata, n = 500L, ...)

## S3 method for class 'adlaplace_fit'
print(x, digits = max(3L, getOption("digits") - 3L), ...)

## S3 method for class 'adlaplace_fit'
summary(object, level = 0.95, ...)

## S3 method for class 'summary.adlaplace_fit'
print(x, digits = max(3L, getOption("digits") - 3L), ...)
```

```
## S3 method for class 'adlaplace_fit'
plot(x, k = 200L, draws = 100L, ...)
```

Arguments

object, x	An adlaplace_fit object.
transform	Logical; report transformed (natural-scale) values.
...	Passed to downstream methods.
parm	Optional subset of parameter labels or indices.
level	Confidence level.
newdata	Data frame of prediction covariates.
n	Number of random-effect draws for predict.
digits	Number of significant digits to print.
k	Number of evaluation points per smooth term in plot.
draws	Number of simulation draws per smooth term in plot.

Details

predict draws random effects from the Laplace (Gaussian) approximation at the mode via [sim_fit](#) and returns the linear predictor on newdata: a nrow(newdata) by n matrix of simulated values (plus fixed-effect contributions from terms whose variables appear in newdata). Without newdata it returns fitted(object).

plot draws each smooth term (terms with more than one knot) on a grid spanning its knots, showing simulated curves from the Laplace approximation with the pointwise mean in red. With no smooth terms it plots the random-effect modes.

apply_theta_log	<i>Apply log transform to selected theta columns</i>
-----------------	--

Description

When active is TRUE, logs values in cols for rows where theta_info\$transform is TRUE. Missing or NA transform entries default to TRUE.

Usage

```
apply_theta_log(theta_info, cols = "init", active = TRUE)
```

Arguments

theta_info	Data frame from info\$theta (see data_setup).
cols	Character vector of column names to transform (e.g. "init").
active	Logical; when FALSE, returns theta_info unchanged.

Value

A copy of `theta_info` with selected entries logged.

binomial-class	<i>Binomial observation term</i>
----------------	----------------------------------

Description

Model term for the observation-level binomial log density with a logit link, registered as `binomial_obs`. For each observation with linear predictor $\eta = X\beta + A\gamma$, trial count N , and success count y , the log density contribution is $y * \eta - N * \log(1 + \exp(\eta))$ (up to a data-only constant). When `size` is omitted, $N = 1$ (Bernoulli). There are no observation-level hyperparameters.

Creates a response term wired to the `binomial_obs` AD shard. A Bernoulli likelihood (`size` omitted) has no observation-level hyperparameters. With `size`, trial counts are read from that column of data and stored as observation weights.

When called with no response variable this function falls back to `stats::binomial()`, so `glm(..., family = binomial)` and `MASS::glmPQL(..., family = binomial)` keep working with **adlaplace** attached.

Usage

```
binomial(x, size = NULL, link = "logit")
```

Arguments

<code>x</code>	Outcome variable name (counts of successes; 0/1 when Bernoulli).
<code>size</code>	Optional column name for the number of trials per observation.
<code>link</code>	Link passed to <code>stats::binomial()</code> in the fallback case.
<code>term</code>	A model term object (S4 methods).
<code>data</code>	A data frame containing the variables (S4 methods).

Details

Binomial observation model term

Value

A binomial object (or a `stats::family` when called with no response variable).

Slots (inherited from model)

`ad_fun` Character scalar "binomial_obs".
`ad_kind` Character scalar "observations".

Slots

`size` Optional column name for per-observation trial counts.

by_group	<i>By-Group Classes and Functions</i>
----------	---------------------------------------

Description

Functions for creating and managing by_group objects for hierarchical model grouping information.

Usage

```
by_group(term, data = NULL, levels = NULL)
```

Arguments

term	Character, the grouping variable name
data	A data frame. If provided and levels is missing, unique values are extracted from data[[term]]
levels	Character vector of unique levels for the grouping variable. If missing and data is provided, unique values are extracted from data[[term]]

Value

A by_group object

Examples

```
# Create with explicit levels
bg <- by_group(term = "group", levels = c("A", "B", "C"))

# Create from data
dat <- data.frame(group = c("A", "B", "A", "C"))
bg <- by_group(term = "group", data = dat)
```

by_group-class	<i>By-Group Class</i>
----------------	-----------------------

Description

A class for grouping information in hierarchical model terms. Contains the grouping variable name and its levels/labels.

Slots

term	Character vector of length 1, the grouping variable name
levels	Character vector of unique levels for the grouping variable
labels	Character vector of formatted labels for the levels

c.ad_fun_ptr	<i>Combine ad_fun_ptr handles</i>
--------------	-----------------------------------

Description

Moves AD shards from one or more `ad_fun_ptr` objects into a single handle. Inputs are cleared after the merge (ownership moves).

Usage

```
## S3 method for class 'ad_fun_ptr'
c(x, ...)
```

Arguments

x	An <code>ad_fun_ptr</code> .
...	Additional <code>ad_fun_ptr</code> objects.

Value

A combined `ad_fun_ptr`.

See Also

[clone_ad_fun_ptr](#), [ad_fun](#)

clone_ad_fun_ptr	<i>Deep copy of an ad_fun_ptr handle</i>
------------------	--

Description

Returns a new `ad_fun_ptr` with independent C++ state (CppAD tapes and sparsity patterns). Hessian templates are not copied; call [ad_fun](#) on the clone to attach maps. The source handle is left unchanged, unlike [c.ad_fun_ptr](#) which moves shards and clears the inputs.

Usage

```
clone_ad_fun_ptr(x)
```

Arguments

x	An <code>ad_fun_ptr</code> object.
---	------------------------------------

Value

A new `ad_fun_ptr`.

collect_terms	<i>Parse Model Terms from Formula</i>
---------------	---------------------------------------

Description

Parses a formula and creates model terms by evaluating constructor calls in the formula environment (so `pkg::term(...)` and terms from attached packages on the search path work). Falls back to the **adlplace** namespace for built-in constructors.

Usage

```
collect_terms(formula, verbose = FALSE)
```

Arguments

formula	Model formula with constructor terms (e.g. <code>iwp(x, ...)</code> , <code>iid(fac)</code> , <code>nbinom(y, ...)</code> on the LHS). Bare symbols such as <code>x</code> are coerced to <code>linear(x)</code> . Unnamed first positional symbols are treated as column names in data; named arguments (e.g. <code>x = sites</code> for spatial terms) are evaluated as objects in the formula environment. The named argument <code>size</code> is also treated as a column name when passed as a symbol (<code>binomial(y, size = N)</code>).
verbose	print extra information

Value

List of model term objects

format_parameters	<i>Format Laplace estimates using parameter metadata</i>
-------------------	--

Description

Maps flat Laplace parameter vectors onto labeled tables from `info` (as stored on `ad_fun@info` or `model_data()$data$info`).

Usage

```
format_parameters(  
  info,  
  full_parameters = NULL,  
  parameters = NULL,  
  gamma = NULL  
)
```

Arguments

info	List with beta, gamma, theta, and parameters data frames.
full_parameters	Numeric vector c(beta, gamma, theta). When supplied, parameters and gamma are ignored.
parameters	Numeric outer vector c(beta, theta); used when full_parameters is NULL.
gamma	Numeric inner vector; used when full_parameters is NULL.

Value

A list with data frames parameters and gamma, or list() when info is missing or incomplete.

fpoly-class	<i>Fixed Polynomial Model Term</i>
-------------	------------------------------------

Description

Creates and manages fixed polynomial model terms.

Usage

```
fpoly(  
  x,  
  p = 2,  
  ref_value = 0,  
  init = .my_beta_init,  
  lower = .my_beta_lower,  
  upper = .my_beta_upper,  
  parscale = .my_beta_parscale  
)  
  
## S4 method for signature 'fpoly'  
design(term, data)  
  
## S4 method for signature 'fpoly'  
beta_info(term, data)
```

Arguments

x	Variable name.
p	Polynomial degree (default: 2).
ref_value	Reference value for the polynomial.
init	Initial values for beta parameters.
lower	Lower bounds for beta parameters.

upper	Upper bounds for beta parameters.
parscale	Parameter scales for optimization.
term	A 'fpoly' term object.
data	A data frame containing the variables used in the term.

Value

A 'fpoly' term object.

A design matrix for the fixed polynomial term, or NULL if p.order is 0.

A data frame containing beta parameter information for the fixed polynomial term.

Functions

- `design(fpoly)`: Design method for fpoly objects
- `beta_info(fpoly)`: Beta info method for fpoly objects

Methods

The following methods are available for 'fpoly' objects:

`design(term, data)` Creates design matrix for fpoly term

`precision(term, data)` Creates precision matrix for fpoly term

`theta_info(term)` Extracts theta parameter information

`beta_info(term, data)` Extracts beta parameter information

`random_info(term, data)` Extracts random effects information

Examples

```
# Create a fixed polynomial term
fpoly_term <- fpoly(x = "temp", p = 3, ref_value = 15)
```

gamm

Example GAMM simulation data

Description

Simulation data for the GAMM vignette example.

Format

A data frame `dat` with 500 rows and columns from `mgcv::gamSim(6, n = 500)`, plus simulated negative-binomial responses used in the GAMM vignette (`y`, `f2`, `mu`, `Z`, etc.).

Source

Generated with `mgcv::gamSim()` and a custom negative-binomial response; see the regeneration comments in `vignettes/gamm.Rmd`.

Examples

```
data("gamm", package = "adlplace")
exists("dat")
nrow(dat)
```

gaussian-class

Gaussian observation term

Description

Model term for the observation-level Gaussian log density registered as `gaussian_obs`. This is the default observation term: a bare response symbol on the left-hand side of a formula (e.g. $y \sim x$) is coerced to `gaussian(y)` by [collect_terms](#).

Creates a response term wired to the `gaussian_obs` AD shard, with a single residual standard deviation parameter (log scale during optimization).

When called with no arguments this function falls back to `stats::gaussian()`, so `glm(..., family = gaussian)` keeps working with **adlplace** attached.

Usage

```
gaussian(
  x,
  init = 1,
  lower = .my_theta_lower,
  upper = .my_theta_upper,
  parscale = .my_theta_parscale
)

## S4 method for signature 'gaussian'
theta_info(term)
```

Arguments

<code>x</code>	Outcome variable name.
<code>init</code>	Initial value for the residual standard deviation.
<code>lower</code>	Lower bound for the residual standard deviation.
<code>upper</code>	Upper bound for the residual standard deviation.
<code>parscale</code>	Parameter scale for optimization.
<code>term</code>	A model term object (S4 methods).
<code>data</code>	A data frame containing the variables (S4 methods).

Details

Gaussian observation model term

Value

A gaussian object (or a `stats::family` when called with no arguments).

Functions

- `theta_info(gaussian)`: Theta info for the residual standard deviation.

Slots (inherited from model)

`ad_fun` Character scalar "gaussian_obs".
`ad_kind` Character scalar "observations".

germany

Germany oral cavity cancer (Besag-York-Mollie example)

Description

Disease-mapping data distributed with **INLA** (Besag, York, and Mollie 1991). Counts y_i in 544 districts follow $y_i \sim \text{Poisson}(E_i e^{\eta_i})$ with a BYM spatial random effect. The bundled ICAR precision `Q_scaled` is the INLA scale.model version with sum-to-zero constraint; `prec` is the `random_mult` precision payload (`Q`, `log_det`, `rank`).

Usage

```
germany
```

Format

A list with:

Y Observed case counts (length 544).

E Expected counts (length 544).

region District index (length 544).

adj District adjacency matrix (`dgCMatrix`, 544 x 544).

Q_scaled Scaled ICAR precision (`dgCMatrix`).

prec List passed to `ad_fun = "random_mult"`: `Q`, `log_det`, and `rank`.

Details

District boundaries are not stored in this object (to avoid a hard dependency on **terra**). They ship as `system.file("extdata", "germany_map.gpkg", package = "adlplace")` and can be loaded with `terra::vect(...)` when **terra** is installed.

Source

INLA Germany data and germany.graph.

References

Besag, J., York, J., and Mollie, A. (1991). Bayesian image restoration, with two applications in spatial statistics. *Annals of the Institute of Statistical Mathematics*, 43(1), 1–20.

Rue, H., Martino, S., and Chopin, N. (2009). Approximate Bayesian inference for latent Gaussian models by using integrated nested Laplace approximations. *Journal of the Royal Statistical Society: Series B*, 71(2), 319–392.

has_openmp	<i>Whether this build was compiled with OpenMP support.</i>
------------	---

Description

Whether this build was compiled with OpenMP support.

Usage

```
has_openmp()
```

Value

TRUE if OpenMP was enabled at compile time, otherwise FALSE.

iid-class	<i>IID Random Effects Term</i>
-----------	--------------------------------

Description

Creates and manages IID (independent and identically distributed) random effects terms.

Usage

```
iid(
  x,
  init = .my_theta_init,
  lower = .my_theta_lower,
  upper = .my_theta_upper,
  parscale = .my_theta_parscale
)

## S4 method for signature 'iid'
design(term, data)
```

```
## S4 method for signature 'iid'
precision(term, data)

## S4 method for signature 'iid'
random_info(term, data)

## S4 method for signature 'iid'
theta_info(term)
```

Arguments

x	Variable name (typically a factor).
init	Initial value for theta parameter.
lower	Lower bound for theta parameter.
upper	Upper bound for theta parameter.
parscale	Parameter scale for optimization.
term	An iid term object
data	A data frame containing the term variable

Value

An ‘iid’ term object (in a named list).

Functions

- `design(iid)`: Creates design matrix for iid term
- `precision(iid)`: Creates precision matrix for iid term
- `random_info(iid)`: Extracts random effects information for iid term
- `theta_info(iid)`: Extracts theta parameter information for iid term

inner_opt

Inner optimization over gamma using trust-region CG (sparse)

Description

Runs the inner optimization over γ using the **trustOptim** sparse trust-region CG solver on a pre-built `ad_fun` handle.

Usage

```
inner_opt(
  parameters,
  gamma,
  ad_fun,
  control = NULL,
  deriv = FALSE,
  verbose = FALSE
)
```

Arguments

parameters	Numeric vector of fixed outer parameters (beta, theta; length Nbeta+Ntheta).
gamma	Numeric vector of starting values for inner parameters (gamma; length Ngamma).
ad_fun	ad_fun S4 object from ad_fun .
control	List of trust-region control parameters (see trustOptim).
deriv	Logical; if TRUE, also return outer gradient/Hessian pieces and Cholesky-based quantities at the inner mode.
verbose	Logical; if TRUE, print thread/shard diagnostics.

Details

Negates tape log-density values in C++ for minimization via **trustOptim**.

Value

A list with log_lik, neg_log_lik, parameters, full_parameters, opt, gradient (list inner/outer), and hessian (list inner/outer/chol_inner/half_log_det; with deriv=TRUE also half_H_inv, H_inv, trace3). Objective and derivatives use the negative log-density convention.

intercept-class	<i>intercept Model Term</i>
-----------------	-----------------------------

Description

Creates and manages intercept model terms for fixed effects.

Usage

```
intercept(
  init = .my_beta_init,
  lower = .my_beta_lower,
  upper = .my_beta_upper,
  parscale = .my_beta_parscale
)
```

```
## S4 method for signature 'intercept'
design(term, data)

## S4 method for signature 'intercept'
beta_info(term, data)
```

Arguments

<code>init</code>	Initial value for beta parameter (default: 0)
<code>lower</code>	Lower bound for beta parameter (default: -Inf)
<code>upper</code>	Upper bound for beta parameter (default: Inf)
<code>parscale</code>	Parameter scale for optimization (default: 1)
<code>term</code>	A intercept term object
<code>data</code>	A data frame containing the term variable

Details

intercept Model Term

Value

A intercept term object

Functions

- `design(intercept)`: Creates design matrix for intercept term
- `beta_info(intercept)`: Extracts beta parameter information for intercept term

Methods

The following methods are available for ‘intercept’ objects:

```
design(term, data)  Creates design matrix for intercept term
precision(term, data)  Creates precision matrix for intercept term
theta_info(term)  Extracts theta parameter information
beta_info(term, data)  Extracts beta parameter information
random_info(term, data)  Extracts random effects information
```

Examples

```
# Create a intercept term
intercept_term <- intercept(init=2)
```

iwp

*Integrated Wiener Process Term Constructor***Description**

Creates an integrated Wiener process (IWP) model term.

Usage

```
iwp(
  x,
  p = 2,
  ref_value = 0,
  knots,
  range = NULL,
  init = .my_theta_init,
  lower = .my_theta_lower,
  upper = .my_theta_upper,
  parscale = .my_theta_parscale,
  boundary_is_random = TRUE,
  include_poly = TRUE
)
```

Arguments

x	Variable name.
p	Order of the integrated Wiener process (default: 2).
ref_value	Reference value for the basis.
knots	Vector of knot locations.
range	Range of the data (optional).
init	Initial values for theta parameters.
lower	Lower bounds for theta parameters.
upper	Upper bounds for theta parameters.
parscale	Parameter scales for optimization.
boundary_is_random	Whether boundary should be treated as random.
include_poly	Whether to include polynomial terms.

Value

A list containing the 'iwp' term object and optionally polynomial terms.

Examples

```
# Example usage:
# knots <- seq(0, 1, length.out = 5)
# iwp_term <- iwp(x = "age", knots = knots)
```

iwp-class

*Integrated Wiener Process Term***Description**

Creates and manages integrated Wiener process (IWP) model terms.

Usage

```
## S4 method for signature 'iwp'
design(term, data)

## S4 method for signature 'iwp'
precision(term, data)

## S4 method for signature 'iwp'
theta_info(term)

## S4 method for signature 'iwp'
random_info(term, data)
```

Arguments

term	An iwp term object
data	A data frame containing the term variable

Functions

- `design(iwp)`: Creates design matrix for IWP term
- `precision(iwp)`: Creates precision matrix for IWP term
- `theta_info(iwp)`: Extracts theta parameter information for IWP term
- `random_info(iwp)`: Extracts random effects information for IWP term

Methods

The following methods are available for ‘iwp’ objects:

```
design(term, data) Creates design matrix for IWP term
precision(term, data) Creates precision matrix for IWP term
theta_info(term) Extracts theta parameter information
beta_info(term, data) Extracts beta parameter information
random_info(term, data) Extracts random effects information
```

laplace_half_H_inv	<i>Extract inner inverse Cholesky factor from Laplace output</i>
--------------------	--

Description

Returns $H^{-1/2}$ for the inner (random-effect) Hessian from the output of `log_lik_laplace`. Accepts the flat `half_H_inv` matrix when `deriv = TRUE`, or reconstructs it from `hessian$chol_inner`.

Usage

```
laplace_half_H_inv(laplace)
```

Arguments

laplace	Output of <code>log_lik_laplace(..., deriv = TRUE)</code> (or <code>inner_opt(..., deriv = TRUE)</code> with the same <code>extra\$hessian</code> layout).
---------	--

Value

A sparse or dense matrix $H^{-1/2}$ with `nrow = Ngamma`.

See Also

[log_lik_laplace](#), [rmvnldl](#)

linear-class	<i>Linear Model Term</i>
--------------	--------------------------

Description

Creates and manages linear model terms for fixed effects.

Usage

```
linear(
  x,
  init = .my_beta_init,
  lower = .my_beta_lower,
  upper = .my_beta_upper,
  parscale = .my_beta_parscale
)

## S4 method for signature 'linear'
design(term, data)

## S4 method for signature 'linear'
beta_info(term, data)
```

Arguments

x	Variable name (character) or symbol in a formula (e.g. linear(x) in model_data()).
init	Initial value for beta parameter (default: 0)
lower	Lower bound for beta parameter (default: -Inf)
upper	Upper bound for beta parameter (default: Inf)
parscale	Parameter scale for optimization (default: 1)
term	A linear term object
data	A data frame containing the term variable

Details

Linear Model Term

Value

A linear term object

Functions

- `design(linear)`: Creates design matrix for linear term
- `beta_info(linear)`: Extracts beta parameter information for linear term

Methods

The following methods are available for 'linear' objects:

`design(term, data)` Creates design matrix for linear term
`precision(term, data)` Creates precision matrix for linear term
`theta_info(term)` Extracts theta parameter information
`beta_info(term, data)` Extracts beta parameter information
`random_info(term, data)` Extracts random effects information

Examples

```
# Create a linear term  
linear_term <- linear(x = "temperature")
```

log_lik_laplace

*Log-likelihood with inner Laplace optimization***Description**

Evaluates the (profiled) log-likelihood for a hierarchical model by solving an inner optimization problem for the latent vector γ (e.g. random effects) given outer parameters x (typically β and θ). The function delegates the inner optimization to `inner_opt()` from the selected backend package (by default **adlaplace**).

Usage

```
log_lik_laplace(
  x,
  config = list(),
  gamma,
  control = list(report.level = 4, report.freq = 1),
  ad_fun,
  data,
  package = c(config[["package"]], "adlaplace")[1L],
  deriv = FALSE
)
```

Arguments

<code>x</code>	Numeric vector of outer parameters (β , then θ). Length must equal <code>num_beta + num_theta</code> from <code>ad_fun@sizes</code> .
<code>config</code>	A list of backend options (verbose, package, etc.). When <code>ad_fun</code> is missing, must also include <code>beta</code> , <code>gamma</code> , and <code>theta</code> so <code>ad_fun()</code> can be built from data.
<code>gamma</code>	Optional numeric vector of starting values for the inner parameter γ . If missing, uses <code>config\$gamma</code> when present, otherwise a zero vector of length <code>ad_fun@sizes["gamma"]</code> .
<code>control</code>	List of control parameters passed <i>as-is</i> to the backend inner optimizer (e.g., <code>report.level</code> , <code>report.freq</code>). See backend documentation (e.g., trustOptim) for supported options.
<code>ad_fun</code>	Optional <code>ad_fun</code> object from ad_fun . This is a single backend handle (no separate inner/outer handles). If missing, it will be constructed automatically using data.
<code>data</code>	Optional data list used to build <code>ad_fun</code> when <code>ad_fun</code> is not supplied.
<code>package</code>	Character scalar naming the backend package to use for <code>ad_fun()</code> and <code>inner_opt()</code> . Defaults to "adlaplace". Other backends must export a compatible <code>ad_fun()</code> builder and <code>inner_opt()</code> .
<code>deriv</code>	Logical scalar. If TRUE, include derivative quantities in the output (gradient, intermediate derivatives).

Details

The default **adlaplace** backend uses a single AD handle. This function passes that handle to `inner_opt(..., ad_fun = ad_fun)`. Parameter block sizes are read from `ad_fun@sizes`; config is not required to contain beta, gamma, or theta when `ad_fun` is supplied.

When `deriv=FALSE`, the return value is the `inner_opt()` list (profile likelihood, nested gradient and hessian, and opt). When `deriv=TRUE`, the same list is augmented with `grad`, `deriv`, and `extra` from the internal profile-derivative helper.

Value

A list. With `deriv=FALSE`, same as `inner_opt`. With `deriv=TRUE`, additionally:

grad Gradient of `neg_log_lik` w.r.t. outer `x`.

deriv Data frame of profile derivative pieces.

extra Intermediate objects (`dU`, `trace3`, Cholesky factors).

Note

The Laplace approximation assumes the inner Hessian is positive definite at the optimum. If `inner_opt` fails to converge or returns a non-invertible Hessian, `log_lik_laplace()` issues a warning or error.

See Also

`ad_fun`, `inner_opt`

mean_mvquad

Posterior mean of random effects via multivariate quadrature

Description

Approximates $E[\gamma \mid y]$ by importance-weighted quadrature over γ , using a Gaussian proposal $N(\hat{\gamma}, H^{-1})$ (Laplace covariance) and correcting with `joint_log_dens`.

Usage

```
mean_mvquad(
  parameters,
  mode,
  cov,
  ad_fun,
  n = 3L,
  type = "nHN",
  ndConstruction = "sparse",
  dec.type = 2L
)
```

Arguments

parameters	Outer (β, θ) at the Laplace fit.
mode	Inner mode $\hat{\gamma}$ (e.g. <code>laplace\$opt\$solution</code>).
cov	H^{-1} for the inner block (e.g. <code>laplace\$extra\$hessian\$H_inv</code>).
ad_fun	<code>ad_fun</code> object for <code>joint_log_dens</code> .
n	Quadrature level passed to mvQuad <code>createNIGrid</code> .
type	<code>mvQuad</code> rule (default "nHN").
ndConstruction	<code>mvQuad</code> grid construction (default "sparse").
dec.type	Cholesky decomposition flag for <code>mvQuad::rescale</code> .

Details

Builds a sparse nHN grid with **mvQuad**, rescales to the Laplace proposal, then forms $\sum_i w_i \gamma_i \exp(\ell(\gamma_i) - \log q(\gamma_i)) / \sum_i w_i \exp(\ell(\gamma_i) - \log q(\gamma_i))$ with log-sum-exp stabilization.

Value

Numeric vector of length `n_gamma`.

See Also

[log_lik_laplace](#), [inner_opt](#), [joint_log_dens](#), [rmvnldl](#)

model-class

Base Model Class

Description

The base S4 class that all specific model term classes inherit from. Subclasses that need hierarchical grouping should include a `by` slot (either character or `by_group` type) and optionally `by_levels` and `by_labels` slots.

Slots

term	Character vector of length 1, the term name
label	Character scalar term label reused across metadata builders.
formula	Formula object for the term
knots	Numeric vector of knot locations (for spline terms)
ref_value	Numeric reference value for centering
p.order	Integer polynomial order
init	Numeric vector of initial values
lower	Numeric vector of lower bounds
upper	Numeric vector of upper bounds

`parscale` Numeric vector of parameter scales for optimization
`type` Factor indicating term type ("fixed", "random", or "response")
`ad_fun` Registered AD density name for `ad_fun_ptr()`, or NA when the term does not map to a density shard.
`ad_kind` Shard kind passed to `ad_fun_ptr()` ("observations", "parameters", "random", or NA).
`package` Package whose `.so` records AD tapes for this term (default "adlaplace" for built-in densities).

model-generics	<i>Model Term Generics</i>
----------------	----------------------------

Description

Generic functions for extracting model matrices and parameter information from model terms.

When non-NULL, `model_data` emits an extra `ad_kind = "parameters"` shard (e.g. FEM log-determinant) alongside the random shard. Default is NULL (no companion).

Usage

```

design(term, data)

precision(term, data)

theta_info(term)

beta_info(term, data)

random_info(term, data)

elgm_matrix(term, data)

## S4 method for signature 'model'
design(term, data)

## S4 method for signature 'model'
precision(term, data)

## S4 method for signature 'model'
theta_info(term)

## S4 method for signature 'model'
beta_info(term, data)

## S4 method for signature 'model'

```

```

random_info(term, data)

extra_ad_fun(term)

## S4 method for signature 'model'
extra_ad_fun(term)

```

Arguments

term	A model term object.
data	A data frame containing the variables.

Value

Method-specific return values.
 Character density name, or NULL.

model_data	<i>Assemble model terms and per-shard ad_data objects</i>
------------	---

Description

Parses a formula, builds design matrices, and returns formula terms plus ad_data shards for the observation density, parameter densities, and each random-effect term.

Usage

```
model_data(formula, data, verbose = FALSE, na_omit = TRUE)
```

Arguments

formula	Model formula with constructor terms (e.g. <code>iwp()</code> , <code>iid()</code>).
data	Data frame containing variables referenced in formula.
verbose	Print extra information while parsing terms.
na_omit	When TRUE (default), drop rows with NA in covariates, outcome, stratification (by), or random-slope mult variables; impute remaining outcome NAs as zero; and sort by observation-term by columns when present.

Value

A list with:

terms Named list of term objects from [collect_terms](#).

data Output of [data_setup](#), including `elgm_matrix` when an observation term defines an `elgm_matrix` method.

observations Named list of observation ad_data shards.

parameters Named list of parameter ad_data shards.

random Named list of random-effect ad_data shards.

Examples

```
## Not run:
md <- model_data(
  y ~ x1 + iwp(x2, p = 2, knots = seq(0, 1, len = 11)),
  data = dat
)

## End(Not run)
```

nbinom-class

*Negative binomial observation term***Description**

Model term for the observation-level negative binomial log density registered as nbinom_obs.
Creates a response term wired to the nbinom_obs AD shard.

Usage

```
nbinom(
  x,
  init = .my_theta_init,
  lower = .my_theta_lower,
  upper = .my_theta_upper,
  parscale = .my_theta_parscale
)

## S4 method for signature 'nbinom'
theta_info(term)
```

Arguments

x	Outcome variable name.
init	Initial value for the overdispersion parameter (log-SD scale).
lower	Lower bound for the overdispersion parameter.
upper	Upper bound for the overdispersion parameter.
parscale	Parameter scale for optimization.
term	A model term object (S4 methods).
data	A data frame containing the variables (S4 methods).

Details

Negative binomial observation model term

Value

A nbinom object.

Functions

- `theta_info(nbinom)`: Theta info for the overdispersion parameter.

Slots (inherited from model)

`ad_fun` Character scalar "nbinom_obs".

`ad_kind` Character scalar "observations".

outer_optim_wrappers *Outer objective and gradient wrappers*

Description

Convenience wrappers around [log_lik_laplace](#) for use with outer optimizers that expect `fn / gr` callbacks. Both functions solve (or warm-start) the inner problem over `gamma` via [log_lik_laplace](#) and update a mutable cache environment with the latest inner solution.

Usage

```
outer_fn(x, config, cache, ad_fun, control_inner = list(), ...)
```

```
outer_gr(x, config, cache, ad_fun, control_inner = list(), ...)
```

Arguments

<code>x</code>	Numeric outer parameter vector <code>c(beta, theta)</code> .
<code>config</code>	Configuration list passed to log_lik_laplace .
<code>cache</code>	An environment used to warm-start and store the inner <code>gamma</code> solution. If <code>cache\$gamma</code> is missing or the wrong length, it is initialized from <code>config\$gamma</code> (if present) or zeros of length <code>ad_fun@sizes["gamma"]</code> . Both functions update <code>cache\$gamma</code> after each evaluation.
<code>ad_fun</code>	<code>ad_fun</code> object from ad_fun .
<code>control_inner</code>	A list of control options forwarded to the <code>control</code> argument of log_lik_laplace for the inner optimization.
<code>...</code>	Additional arguments forwarded to log_lik_laplace (for example data or package).

Details

`outer_fn()` Returns the scalar objective negative log likelihood.

`outer_gr()` Returns the gradient.

Value

- outer_fn: a numeric scalar neg_log_lik.
- outer_gr: gradient of neg_log_lik (same sign as outer_fn).

See Also

[log_lik_laplace](#)

Examples

```
## Not run:
cache <- new.env(parent = emptyenv())
cache$gamma <- rep(0, nrow(data$ATp))
ad_fun <- ad_fun(data, config)

val <- outer_fn(x = x0, data = data, config = config, cache = cache, ad_fun = ad_fun)
gr <- outer_gr(x = x0, data = data, config = config, cache = cache, ad_fun = ad_fun)

## End(Not run)
```

rmvnldl

Simulate from a multivariate normal using LDL of the precision matrix

Description

Draws n samples from $N(\mu, H^{-1})$ where H is the precision matrix, given its LDL factorization (unit lower L1, diagonal D, and permutation perm).

Usage

```
rmvnldl(n, mean = 0, chol_prec)
```

Arguments

n	Number of samples.
mean	Mean vector μ (length length(D)).
chol_prec	Either a CHMfactor, Cholesky, or dCHMsimpl from <code>Matrix::Cholesky(..., LDL = TRUE)</code> , or a list like <code>res\$hessian\$chol_inner</code> with components L1, D, and perm.

Details

Uses $H^{-1/2} = P^T L^{-1} D^{-1/2}$ (same as `reformat_chol()` in `log_lik_deriv`), and `rnorm` only: $x = \mu + H^{-1/2}z, z \sim N(0, I)$.

Value

An n-by-p matrix; each row is one draw.

See Also

[Cholesky](#), [log_lik_laplace](#)

rpoly-class	<i>Random Polynomial Model Term</i>
-------------	-------------------------------------

Description

Creates and manages random polynomial model terms.

Usage

```
rpoly(x, p = 2, ref_value = 0, sd = Inf)

## S4 method for signature 'rpoly'
design(term, data)

## S4 method for signature 'rpoly'
precision(term, data)

## S4 method for signature 'rpoly'
random_info(term, data)
```

Arguments

- x Variable name.
- p Polynomial degree (default: 2).
- ref_value Reference value for the polynomial.
- sd Standard deviation for random effects.
- term A ‘rpoly’ term object.
- data A data frame containing the variables used in the term.

Value

- A ‘rpoly’ term object.
- A design matrix for the random polynomial term.
- A precision matrix for the random polynomial term.
- A data frame containing random effects information for the random polynomial term.

Functions

- `design(rpoly)`: Design method for `rpoly` objects
- `precision(rpoly)`: Precision method for `rpoly` objects
- `random_info(rpoly)`: Random info method for `rpoly` objects

Methods

The following methods are available for 'rpoly' objects:

`design(term, data)` Creates design matrix for `rpoly` term
`precision(term, data)` Creates precision matrix for `rpoly` term
`theta_info(term)` Extracts theta parameter information
`beta_info(term, data)` Extracts beta parameter information
`random_info(term, data)` Extracts random effects information

sim_fit

Simulate linear predictors on a new covariate grid

Description

Draws random effects from the Laplace approximation and evaluates the linear predictor on `x` by summing fixed (β) and random (γ) contributions from each model term.

Usage

```
sim_fit(x, data, fit, n = 500L)
```

Arguments

<code>x</code>	Data frame of prediction covariates (same variables as the fitted model). Only terms whose variables appear in <code>x</code> contribute.
<code>data</code>	Object returned by <code>model_data</code> .
<code>fit</code>	Result from <code>log_lik_laplace</code> at the fitted outer parameters.
<code>n</code>	Number of draws for random-effect simulation.

Value

The combined linear predictor: a `nrow(x)` by `n` matrix of simulated values (fixed-effect contributions are constant across columns).

sizes	<i>Parameter-block row counts for an ad_data</i>
-------	--

Description

Parameter-block row counts for an `ad_data`

Usage

```
sizes(x, ...)  
  
## S4 method for signature 'ad_data'  
sizes(x)
```

Arguments

<code>x</code>	An <code>ad_data</code> object.
<code>...</code>	Unused; for S4 generic compatibility.

Value

Named list with `n_beta`, `n_gamma`, and `n_theta`.

Index

- * **datasets**
 - .type_factor_levels, 3
 - gamm, 19
 - germany, 21
 - .type_factor_levels, 3
- ad_data, 3
- ad_data-class, 5
- ad_fun, 5, 7–9, 16, 23, 24, 30, 31, 36
- ad_fun, ad_data-method (ad_fun), 5
- ad_fun, ad_fun_ptr-method (ad_fun), 5
- ad_fun, list-method (ad_fun), 5
- ad_fun-class, 6
- ad_fun_ptr, 7, 8
- ad_fun_ptr-class, 8
- ad_shards, 8, 9
- adlaplace, 9, 12
- adlaplace_cpp, 11
- adlaplace_fit-methods, 12
- apply_theta_log, 13
- beta_info (model-generics), 33
- beta_info, fpoly-method (fpoly-class), 18
- beta_info, intercept-method (intercept-class), 24
- beta_info, linear-method (linear-class), 28
- beta_info, model-method (model-generics), 33
- binomial (binomial-class), 14
- binomial-class, 14
- by_group, 15
- by_group-class, 15
- c.ad_fun_ptr, 8, 16, 16
- Cholesky, 38
- clone_ad_fun_ptr, 8, 16, 16
- coef.adlaplace_fit (adlaplace_fit-methods), 12
- collect_terms, 9, 17, 20, 34
- confint.adlaplace_fit, 10
- confint.adlaplace_fit (adlaplace_fit-methods), 12
- dat (gamm), 19
- data_setup, 13, 34
- data_setup (ad_data), 3
- design (model-generics), 33
- design, fpoly-method (fpoly-class), 18
- design, iid-method (iid-class), 22
- design, intercept-method (intercept-class), 24
- design, iwp-method (iwp-class), 27
- design, linear-method (linear-class), 28
- design, model-method (model-generics), 33
- design, rpoly-method (rpoly-class), 38
- elgm_matrix (model-generics), 33
- environment, 36
- extra_ad_fun (model-generics), 33
- extra_ad_fun, model-method (model-generics), 33
- fitted.adlaplace_fit (adlaplace_fit-methods), 12
- format_parameters, 17
- fpoly (fpoly-class), 18
- fpoly-class, 18
- fun_obj_fdfh (adlaplace_cpp), 11
- gamm, 19
- gaussian (gaussian-class), 20
- gaussian-class, 20
- germany, 21
- grad (adlaplace_cpp), 11
- has_openmp, 22
- hessian (adlaplace_cpp), 11
- iid (iid-class), 22
- iid-class, 22

inner_opt, [23](#), [31](#), [32](#)
 intercept (intercept-class), [24](#)
 intercept-class, [24](#)
 iwp, [26](#)
 iwp-class, [27](#)

 joint_log_dens, [31](#), [32](#)
 joint_log_dens (adlaplace_cpp), [11](#)

 laplace_half_H_inv, [28](#)
 linear (linear-class), [28](#)
 linear-class, [28](#)
 log_lik_laplace, [10](#), [28](#), [30](#), [32](#), [36–39](#)
 logLik.adlaplace_fit
 (adlaplace_fit-methods), [12](#)

 mean_mvquad, [31](#)
 model-class, [32](#)
 model-generics, [33](#)
 model_data, [9](#), [10](#), [33](#), [34](#), [39](#)

 nbinom (nbinom-class), [35](#)
 nbinom-class, [35](#)
 nobs.adlaplace_fit
 (adlaplace_fit-methods), [12](#)

 optim, [9](#), [10](#)
 outer_fn, [9](#)
 outer_fn (outer_optim_wrappers), [36](#)
 outer_gr, [9](#)
 outer_gr (outer_optim_wrappers), [36](#)
 outer_optim_wrappers, [36](#)

 plot.adlaplace_fit
 (adlaplace_fit-methods), [12](#)
 precision (model-generics), [33](#)
 precision, iid-method (iid-class), [22](#)
 precision, iwp-method (iwp-class), [27](#)
 precision, model-method
 (model-generics), [33](#)
 precision, rpoly-method (rpoly-class), [38](#)
 predict.adlaplace_fit, [10](#)
 predict.adlaplace_fit
 (adlaplace_fit-methods), [12](#)
 print.adlaplace_fit
 (adlaplace_fit-methods), [12](#)
 print.summary.adlaplace_fit
 (adlaplace_fit-methods), [12](#)

 random_info (model-generics), [33](#)
 random_info, iid-method (iid-class), [22](#)
 random_info, iwp-method (iwp-class), [27](#)
 random_info, model-method
 (model-generics), [33](#)
 random_info, rpoly-method (rpoly-class),
 [38](#)
 rmvnlDL, [28](#), [32](#), [37](#)
 rpoly (rpoly-class), [38](#)
 rpoly-class, [38](#)

 sim_fit, [13](#), [39](#)
 sizes, [40](#)
 sizes, ad_data-method (sizes), [40](#)
 summary.adlaplace_fit, [10](#)
 summary.adlaplace_fit
 (adlaplace_fit-methods), [12](#)
 svd, [8](#)

 theta_info (model-generics), [33](#)
 theta_info, gaussian-method
 (gaussian-class), [20](#)
 theta_info, iid-method (iid-class), [22](#)
 theta_info, iwp-method (iwp-class), [27](#)
 theta_info, model-method
 (model-generics), [33](#)
 theta_info, nbinom-method
 (nbinom-class), [35](#)
 trace_hinv_t (adlaplace_cpp), [11](#)

 vcov.adlaplace_fit, [10](#)
 vcov.adlaplace_fit
 (adlaplace_fit-methods), [12](#)