Package: postcard (via r-universe)

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Type Package

Title Estimating Marginal Effects with Prognostic Covariate Adjustment

Version 1.1.0

Description Conduct power analyses and inference of marginal effects. Uses plug-in estimation and influence functions to perform robust inference, optionally leveraging historical data to increase precision with prognostic covariate adjustment. The methods are described in Højbjerre-Frandsen et al. (2025) <doi:10.48550/arXiv.2503.22284>.

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Contents

`default_learners` *Creates a list of learners*

Description

This function creates a list of learners compatible with the `learners` argument of fit_best_learner, which is used as the default argument.

Usage

```
default_learners()
```

Value

a named `list` of learners, where each element consists of a

- `model`: A `parsnip` model specification
- `grid`: A `data.frame` with columns corresponding to tuning parameters

Examples

```
default_learners()
```

`fit_best_learner` *Find the best learner in terms of RMSE among specified learners using cross validation*

Description

Find the best learner in terms of RMSE among specified learners using cross validation

Usage

```
fit_best_learner(
  preproc,
  data,
  cv_folds = 5,
  learners = default_learners(),
  verbose = options::opt("verbose")
)
```

Arguments

`preproc`	A list (preferably named) with preprocessing objects: formulas, recipes, or `workflows::workflow_variables()`. Passed to `workflowsets::workflow_set()`.
`data`	A data frame.
`cv_folds`	a `numeric` with the number of cross-validation folds used when fitting and evaluating models
`learners`	a `list` (preferably named) containing named lists of elements `model` and optionally `grid`. The `model` element should be a `parsnip` model specification, which is passed to workflowsets::workflow_set as the `model` argument, while the `grid` element is passed as the `grid` argument of workflowsets::option_add
`verbose`	`numeric` verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution (Defaults to 2, overwritable using option 'postcard.verbose' or environment variable 'R_POSTCARD_VERBOSE')

Details

Ensure data compatibility with the learners.

Value

a trained `workflow`

See Also

See `rctglm_with_prognosticscore()` for a function that utilises this function to perform prognostic covariate adjustment.

Examples

```
# Generate some synthetic 2-armed RCT data along with historical controls
n <- 100
dat_rct <- glm_data(
  Y ~ 1+2*x1+3*a,
  x1 = rnorm(n, 2),
  a = rbinom (n, 1, .5),
  family = gaussian()
)
dat_hist <- glm_data(
  Y ~ 1+2*x1,
  x1 = rnorm(n, 2),
  family = gaussian()
)

# Fit a learner to the historical control data
learners <- list(
  mars = list(
    model = parsnip::set_engine(
      parsnip::mars(
        mode = "regression", prod_degree = 3
      ),
      "earth"
    )
  )
)
fit <- fit_best_learner(
  preproc = list(mod = Y ~ .),
  data = dat_hist,
  learners = learners
)

# Use it fx. to predict the "control outcome" in the 2-armed RCT
predict(fit, new_data = dat_rct)
```

`glm_data` *Generate data simulated from a GLM*

Description

Provide a formula, variables and a family to generate a linear predictor using the formula and provided variables before using the inverse link of the family to generate the GLM modelled mean, mu, which is then used to simulate the response with this mean from the generating function according to the chosen family.

Usage

```
glm_data(formula, ..., family = gaussian(), family_args = NULL)
```

Arguments

formula	an object of class "formula" (or one that can be coerced to that class): a symbolic description of the model to be fitted. The details of model specification are given under 'Details' in the glm documentation.
...	a data.frame with columns corresponding to variables used in formula, a named list of those variables, or individually provided named arguments of variables
family	the family of the response. this can be a character string naming a family function, a family function or the result of a call to a family function
family_args	a named list with values of arguments passed to family relevant r<family_name> function for simulating the data

Value

a data.frame

Examples

```
# Generate a gaussian response from a single covariate
glm_data(Y ~ 1+2*x1,
                x1 = rnorm(10))

# Generate a gaussian response from a single covariate with non-linear
# effects. Specify that the response should have standard deviation sqrt(3)
glm_data(Y ~ 1+2*log(x1),
                x1 = runif(10, min = 1, max = 10),
                family_args = list(sd = sqrt(3)))

# Generate a negative binomial response
glm_data(Y ~ 1+2*x1-x2,
                x1 = rnorm(10),
                x2 = rgamma(10, shape = 2),
                family = MASS::negative.binomial(2))

# Provide variables as a list/data.frame and pass a link to the negative.binomial
# function
glm_data(resp ~ 1+2*x1-x2,
                data.frame(
                  x1 = rnorm(10),
                  x2 = rgamma(10, shape = 2)
                ),
                family = MASS::negative.binomial(2),
                family_args = list(link = "identity"))
```

`options` *postcard Options*

Description

Internally used, package-specific options. All options will prioritize R options() values, and fall back to environment variables if undefined. If neither the option nor the environment variable is set, a default value is used.

Arguments

`verbose`	`numeric` verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution (Defaults to 2, overwritable using option 'postcard.verbose' or environment variable 'R_POSTCARD_VERBOSE')

Checking Option Values

Option values specific to `postcard` can be accessed by passing the package name to `env`.

```
options::opts(env = "postcard")

options::opt(x, default, env = "postcard")
```

Options

verbose `numeric` verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution

default: `2`

option: postcard.verbose

envvar: R_POSTCARD_VERBOSE (evaluated if possible, raw string otherwise)

See Also

options getOption Sys.setenv Sys.getenv

power_linear *Power and sample size estimation for linear models*

Description

variance_ancova provides a convenient function for estimating a variance to use for power and sample size approximation.

The power_gs and samplesize_gs functions calculate the Guenther-Schouten power approximation for ANOVA or ANCOVA. The approximation is based in (Guenther WC. Sample Size Formulas for Normal Theory T Tests. The American Statistician. 1981;35(4):243–244) and (Schouten HJA. Sample size formula with a continuous outcome for unequal group sizes and unequal variances. Statistics in Medicine. 1999;18(1):87–91).

The function power_nc calculates the power for ANOVA or ANCOVA based on the non-centrality parameter and the exact t-distributions.

See more details about each function in Details and in sections after Value.

Usage

```
variance_ancova(formula, data, inflation = 1, deflation = 1, ...)

power_gs(variance, ate, n, r = 1, margin = 0, alpha = 0.05, ...)

samplesize_gs(variance, ate, r = 1, margin = 0, power = 0.9, alpha = 0.05, ...)

power_nc(variance, df, ate, n, r = 1, margin = 0, alpha = 0.05, ...)
```

Arguments

formula	an object of class "formula" (or one that can be coerced to that class): a symbolic description used in stats::model.frame() to create a data.frame with response and covariates. This data.frame is used to estimate the R^2, which is then used to find the variance. See more in details.
data	a data frame, list or environment (or object coercible by as.data.frame to a data frame), containing the variables in formula. Neither a matrix nor an array will be accepted.
inflation	a numeric to multiply the marginal variance of the response by. Default is 1 which estimates the variance directly from data. Use values above 1 to obtain a more conservative estimate of the marginal response variance.
deflation	a numeric to multiply the R^2 by. Default is 1 which means the estimate of R^2 is unchanged. Use values below 1 to obtain a more conservative estimate of the coefficient of determination. See details about how R^2 related to the estimation.
...	Not currently used. Here to accomodate implementation of ... for the repeat_power_linear() function.
variance	a numeric variance to use for the approximation. See more details in documentation sections of each power approximating function.

ate	a numeric minimum effect size that we should be able to detect.
n	a numeric with number of participants in total. From this number of participants in the treatment group is $n1 = (r/(1+r))n$ and the control group is $n0 = (1/(1+r))n$.
r	a numeric allocation ratio $r = n1/n0$. For one-to-one randomisation r=1.
margin	a numeric superiority margin (for non-inferiority margin, a negative value can be provided).
alpha	a numeric significance level. The critical value used for testing corresponds to using the significance level for a two-sided test. I.e. we use the quantile $1-\alpha/2$ of the null distribution as the critical value.
power	a numeric giving the desired power when calculating the sample size
df	a numeric degrees of freedom to use in the t-distribution.

Details

This details section provides information about relation between arguments to functions and the formulas described in sections below for each power approximation formula.

Note that all entities that carry the same name as an argument and in the formula will not be mentioned below, as they are obviously linked (n, r, alpha)

- ate: $\beta_1 - \beta_0$
- margin: Δ_s
- variance: $\widehat{\sigma}^2(1-\widehat{R}^2)$

Finding the variance **to use for approximation:**
The variance_ancova function estimates $\sigma^2(1-R^2)$ in data and returns it as a numeric that can be passed directly as the variance in power_gs. Corresponds to estimating the power from using an lm with the same formula as specified in variance_ancova.
The user can estimate the variance any way they see fit.

Value

All functions return a numeric. variance_ancova returns a numeric with a variance estimated from data to used for power estimation and sample size estimation. power_xx and samplesize_xx functions return a numeric with the power or sample size approximation.

Guenther-Schouten power approximation

The estimation formula in the case of an ANCOVA model with multiple covariate adjustment is (see description for reference):

$$n = \frac{(1+r)^2}{r}\frac{(z_{1-\alpha/2}+z_{1-\beta})^2\widehat{\sigma}^2(1-\widehat{R}^2)}{(\beta_1-\beta_0-\Delta_s)^2} + \frac{(z_{1-\alpha/2})^2}{2}$$

where $\widehat{R}^2 = \frac{\widehat{\sigma}_{XY}^\top \widehat{\Sigma}_X^{-1}\widehat{\sigma}_{XY}}{\widehat{\sigma}^2}$, we denote by $\widehat{\sigma^2}$ an estimate of the variance of the outcome, $\widehat{\Sigma_X}$ and estimate of the covariance matrix of the covariates, and $\widehat{\sigma_{XY}}$ a p-dimensional column vector consisting of an estimate of the covariance between the outcome variable and each covariate. In the univariate case R^2 is replaced by ρ^2

Power approximation using non-centrality parameter

The prospective power estimations are based on (Kieser M. Methods and Applications of Sample Size Calculation and Recalculation in Clinical Trials. Springer; 2020). The ANOVA power is calculated based on the non-centrality parameter given as

$$nc = \sqrt{\frac{r}{(1+r)^2} \cdot n} \cdot \frac{\beta_1 - \beta_0 - \Delta_s}{\sigma},$$

where we denote by σ^2 the variance of the outcome, such that the power can be estimated as

$$1 - \beta = 1 - F_{t,n-2,nc}\left(F^{-1}_{t,n-2,0}(1-\alpha/2)\right).$$

The power of ANCOVA with univariate covariate adjustment and no interaction is calculated based on the non-centrality parameter given as

$$nc = \sqrt{\frac{rn}{(1+r)^2}} \frac{\beta_1 - \beta_0 - \Delta_s}{\sigma\sqrt{1-\rho^2}},$$

such that the power can be estimated as

$$1 - \beta = 1 - F_{t,n-3,nc}\left(F^{-1}_{t,n-3,0}(1-\alpha/2)\right).$$

The power of ANCOVA with either univariate covariate adjustment and interaction or multiple covariate adjustement with or without interaction is calculated based on the non-centrality parameter given as

$$nc = \frac{\beta_1 - \beta_0 - \Delta_s}{\sqrt{\left(\frac{1}{n_1} + \frac{1}{n_0} + X_d^\top \left((n-2)\Sigma_X\right)^{-1} X_d\right)\sigma^2\left(1-\widehat{R}^2\right)}}.$$

where $X_d = \left(\overline{X}^1_1 - \overline{X}^1_0, \ldots, \overline{X}^p_1 - \overline{X}^p_0\right)^\top$, $\widehat{R}^2 = \frac{\widehat{\sigma}^\top_{XY}\widehat{\Sigma}^{-1}_X\widehat{\sigma}_{XY}}{\widehat{\sigma}^2}$, we denote by $\widehat{\sigma^2}$ an estimate of the variance of the outcome, $\widehat{\Sigma_X}$ and estimate of the covariance matrix of the covariates, and $\widehat{\sigma_{XY}}$ a p-dimensional column vector consisting of an estimate of the covariance between the outcome variable and each covariate. Since we are in the case of randomized trials the expected difference between the covariate values between the to groups is 0. Furthermore, the elements of Σ^{-1}_X will be small, unless the variances are close to 0, or the covariates exhibit strong linear dependencies, so that the correlations are close to 1. These scenarios are excluded since they could lead to potentially serious problems regarding inference either way. These arguments are used by Zimmermann et. al (Zimmermann G, Kieser M, Bathke AC. Sample Size Calculation and Blinded Recalculation for Analysis of Covariance Models with Multiple Random Covariates. Journal of Biopharmaceutical Statistics. 2020;30(1):143–159.) to approximate the non-centrality parameter as in the univariate case where ρ^2 is replaced by R^2.

Then the power for ANCOVA with d degrees of freedom can be estimated as

$$1 - \beta = 1 - F_{t,d,nc}\left(F^{-1}_{t,d,0),0}(1-\alpha/2)\right).$$

See Also

See power_marginaleffect for a power approximation function that works for a larger class of models.

Examples

```
# Generate a data set to use as an example
n_train <- 1e3
n_test <- 100

dat_gaus <- glm_data(Y ~ 1+2*X1-X2+3*A,
                X1 = rnorm(n_train),
                X2 = rgamma(n_train, shape = 2),
                A = rbinom(n_train, size = 1, prob = 0.5),
                family = gaussian())

# Approximate the power using no adjustment covariates
va_nocov <- var(dat_gaus$Y)
power_gs(n = n_test, variance = va_nocov, ate = 1)

# Approximate the power with a model adjusting for both variables in the
# data generating process

## First estimate the variance sigma^2 * (1-R^2)
va_cov <- variance_ancova(Y ~ X1 + X2 + A, dat_gaus)
## Then estimate the power using this variance
power_gs(n = n_test, variance = va_cov, ate = 1.8, margin = 1, r = 2)

# Approximate the sample size needed to obtain 90% power with same model as
# above
samplesize_gs(
  variance = va_cov, ate = 1.8, power = 0.9, margin = 1, r = 2
)

# No adjustment covariates
power_nc(n = n_test, variance = va_nocov, df = 199, ate = 1)
# Adjusting for all covariates in data generating process
power_nc(n = n_test, variance = va_cov, df = 196, ate = 1.8, margin = 1, r = 2)
```

---

power_marginaleffect *Power approximation for estimating marginal effects in GLMs*

Description

The functions implements the algorithm for power estimation described in Powering RCTs for marginal effects with GLMs using prognostic score adjustment by Højbjerre-Frandsen et. al (2025). See a bit more context in details and all details in reference.

Usage

```
power_marginaleffect(
  response,
  predictions,
  target_effect,
  exposure_prob,
  var1 = NULL,
  kappa1_squared = NULL,
  estimand_fun = "ate",
  estimand_fun_deriv0 = NULL,
  estimand_fun_deriv1 = NULL,
  inv_estimand_fun = NULL,
  margin = estimand_fun(1, 1),
  alpha = 0.05,
  tolerance = 0.001,
  verbose = options::opt("verbose"),
  ...
)
```

Arguments

`response`	a vector of mode `numeric` with the response variable from comparator participants
`predictions`	a vector of mode `numeric` with predictions of the response
`target_effect`	a `numeric` minimum effect size that we should be able to detect. See more in details.
`exposure_prob`	a `numeric` with the probability of being in "group 1" (rather than group 0). See more in details.
`var1`	a `numeric` variance of the potential outcome corresponding to group 1, or a `function` with a single argument meant to obtain `var1` as a transformation of the variance of the potential outcome corresponding to group 0. See more in details.
`kappa1_squared`	a `numeric` mean-squared error from predicting potential outcome corresponding to group 1, or a `function` with a single arguments meant to obtain `kappa1_squared` as a transformation of the MSE in group 0. See more in details.
`estimand_fun`	a `function` with arguments `psi1` and `psi0` specifying the estimand. Alternative, specify "ate" or "rate_ratio" as a `character` to use one of the default estimand functions. See more details in the "Estimand" section of rctglm.
`estimand_fun_deriv0`	a `function` specifying the derivative of `estimand_fun` wrt. `psi0`. As a default the algorithm will use symbolic differentiation to automatically find the derivative from `estimand_fun`
`estimand_fun_deriv1`	a `function` specifying the derivative of `estimand_fun` wrt. `psi1`. As a default the algorithm will use symbolic differentiation to automatically find the derivative from `estimand_fun`

`inv_estimand_fun`
: (optional) a `function` with arguments `psi0` and `target_effect`, so `estimand_fun(psi1 = y, psi0 = x) = z` and `inv_estimand_fun(psi0 = x, target_effect = z) = y` for all x, y, z. If left as `NULL`, the inverse will be found numerically.

`margin`
: a `numeric` superiority margin. As a default, the `estimand_fun` is evaluated with the same counterfactual means `psi1` and `psi0`, corresponding to a superiority margin assuming no difference (fx. 0 for ATE and 1 for rate ratio).

`alpha`
: a `numeric` significance level. The critical value used for testing corresponds to using the significance level for a two-sided test. I.e. we use the quantile $1 - \alpha/2$ of the null distribution as the critical value.

`tolerance`
: passed to all.equal when comparing calculated `target_effect` from derivations and given `target_effect` during numeric derivation of `inv_estimand_fun`. Thus only relevant when `inv_estimand_fun` is `NULL`.

`verbose`
: `numeric` verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution (Defaults to 2, overwritable using option 'postcard.verbose' or environment variable 'R_POSTCARD_VERBOSE')

`...`
: arguments passed to stats::uniroot, which is used to find the inverse of `estimand_fun`, when `inv_estimand_fun` is `NULL`.

Details

The reference in the description shows in its "Prospective power" section a derivation of a variance bound

$$v_\infty^{\uparrow 2} = r_0'^2\sigma_0^2 + r_1'^2\sigma_1^2 + \pi_0\pi_1\left(\frac{|r_0'|\kappa_0}{\pi_0} + \frac{|r_1'|\kappa_1}{\pi_1}\right)^2$$

where r_a' is the derivative of the `estimand_fun` with respect to Ψ_a, σ_a^2 is the variance of the potential outcome corresponding to group a, π_a is the probability of being assigned to group a, and κ_a is the expected mean-squared error when predicting the potential outcome corresponding to group a.

The variance bound is then used for calculating a lower bound of the power using the distributions corresponding to the null and alternative hypotheses $\mathcal{H}_0 : \hat{\Psi} \sim F_0 = \mathcal{N}(\Delta, v_\infty^{\uparrow 2}/n)$ and $\mathcal{H}_1 : \hat{\Psi} \sim F_1 = \mathcal{N}(\Psi, v_\infty^{\uparrow 2}/n)$. See more details in the reference.

Relating arguments to formulas:

- `response`: Used to obtain both σ_0^2 (by taking the sample variance of the response) and κ_0.
- `predictions`: Used when calculating the MSE κ_0.
- `var1`: σ_1^2. As a default, chosen to be the same as σ_0^2. Can specify differently through this argument fx. by
  - Inflating or deflating the value of σ_0^2 by a scalar according to prior beliefs. Fx. specify `var1 = function(x) 1.2 * x` to inflate σ_0^2 by 1.2.
  - If historical data is available for group 1, an estimate of the variance from that data can be provided here.
- `kappa1_squared`: κ_1. Same as for `var1`, default is to assume the same value as `kappa0_squared`, which is found by using the `response` and `predictions` vectors. Adjust the value according to prior beliefs if relevant.
- `target_effect`: Ψ.
- `exposure_prob`: π_1

Value

a numeric with the estimated power.

See Also

See power_linear for power approximation functionalities specific to linear models.

Examples

```
# Generate a data set to use as an example
n <- 100
exposure_prob <- .5

dat_gaus <- glm_data(Y ~ 1+2*X1-X2+3*A+1.6*A*X2,
                X1 = rnorm(n),
                X2 = rgamma(n, shape = 2),
                A = rbinom(n, size = 1, prob = exposure_prob),
                family = gaussian())

# ------------------------------------------------------------------------
# Obtain out-of-sample (OOS) prediction using glm model
# ------------------------------------------------------------------------
gaus1 <- dat_gaus[1:(n/2), ]
gaus2 <- dat_gaus[(n/2+1):n, ]

glm1 <- glm(Y ~ X1 + X2 + A, data = gaus1)
glm2 <- glm(Y ~ X1 + X2 + A, data = gaus2)
preds_glm1 <- predict(glm2, newdata = gaus1, type = "response")
preds_glm2 <- predict(glm1, newdata = gaus2, type = "response")
preds_glm <- c(preds_glm1, preds_glm2)

# Obtain power
power_marginaleffect(
  response = dat_gaus$Y,
  predictions = preds_glm,
  target_effect = 2,
  exposure_prob = exposure_prob
)

# ------------------------------------------------------------------------
# Get OOS predictions using discrete super learner and adjust variance
# ------------------------------------------------------------------------
learners <- list(
  mars = list(
    model = parsnip::set_engine(
      parsnip::mars(
        mode = "regression", prod_degree = 3
      ),
      "earth"
    )
 ),
    lm = list(
```

```
      model = parsnip::set_engine(
        parsnip::linear_reg(),
        "lm"
      )
    )
)
lrnr1 <- fit_best_learner(preproc = list(mod = Y ~ X1 + X2 + A),
                          data = gaus1,
                          learners = learners)
lrnr2 <- fit_best_learner(preproc = list(mod = Y ~ X1 + X2 + A),
                          data = gaus2,
                          learners = learners)
preds_lrnr1 <- dplyr::pull(predict(lrnr2, new_data = gaus1))
preds_lrnr2 <- dplyr::pull(predict(lrnr1, new_data = gaus2))
preds_lrnr <- c(preds_lrnr1, preds_lrnr2)

# Estimate the power AND adjust the assumed variance in the "unknown"
# group 1 to be 20% larger than in group 0
power_marginaleffect(
  response = dat_gaus$Y,
  predictions = preds_lrnr,
  target_effect = 2,
  exposure_prob = exposure_prob,
  var1 = function(var0) 1.2 * var0
)
```

`prog` *Extract information about the fitted prognostic model*

Description

Extracts the `prognostic_info` list element from an `rctglm_prog` object. See 'Value' at rctglm_with_prognosticscore for more details.

Usage

```
prog(x)

## S3 method for class 'rctglm_prog'
prog(x)
```

Arguments

`x`	an object of class `rctglm_prog` (returned by rctglm_with_prognosticscore)

Value

a list with the structure described of `prognostic_info` in the `Value` section of rctglm_with_prognosticscore.

See Also

The generic rctglm_with_prognosticscore() for which this method works.

Examples

```
# Generate some data
n <- 100
b0 <- 1
b1 <- 1.5
b2 <- 2
W1 <- runif(n, min = 1, max = 10)
exposure_prob <- .5

dat_treat <- glm_data(
  Y ~ b0+b1*log(W1)+b2*A,
  W1 = W1,
  A = rbinom(n, 1, exposure_prob)
)

dat_notreat <- glm_data(
  Y ~ b0+b1*log(W1),
  W1 = W1
)

learners <- list(
  mars = list(
    model = parsnip::set_engine(
      parsnip::mars(
        mode = "regression", prod_degree = 3
      ),
      "earth"
    )
  )
)
ate <- rctglm_with_prognosticscore(
  formula = Y ~ .,
  exposure_indicator = A,
  exposure_prob = exposure_prob,
  data = dat_treat,
  family = gaussian(),
  estimand_fun = "ate",
  data_hist = dat_notreat,
  learners = learners)

prog(ate)
```

---

rctglm *Fit GLM and find any estimand (marginal effect) using plug-in estimation with variance estimation using influence functions*

Description

The procedure uses plug-in-estimation and influence functions to perform robust inference of any specified estimand in the setting of a randomised clinical trial, even in the case of heterogeneous effect of covariates in randomisation groups. See Powering RCTs for marginal effects with GLMs using prognostic score adjustment by Højbjerre-Frandsen et. al (2025) for more details on methodology.

Usage

```
rctglm(
  formula,
  exposure_indicator,
  exposure_prob,
  data,
  family = gaussian,
  estimand_fun = "ate",
  estimand_fun_deriv0 = NULL,
  estimand_fun_deriv1 = NULL,
  cv_variance = FALSE,
  cv_variance_folds = 10,
  verbose = options::opt("verbose"),
  ...
)
```

Arguments

`formula`	an object of class "formula" (or one that can be coerced to that class): a symbolic description of the model to be fitted. The details of model specification are given under 'Details' in the glm documentation.
`exposure_indicator`	(name of) the *binary* variable in `data` that identifies randomisation groups. The variable is required to be binary currently.
`exposure_prob`	a `numeric` with the probability of being in "group 1" (rather than group 0) in groups defined by `exposure_indicator`.
`data`	an optional data frame, list or environment (or object coercible by as.data.frame to a data frame) containing the variables in the model. If not found in data, the variables are taken from environment(formula), typically the environment from which the function is called.
`family`	a description of the error distribution and link function to be used in the model. For `glm` this can be a character string naming a family function, a family function or the result of a call to a family function. For `glm.fit` only the third option is supported. (See `family` for details of family functions.)
`estimand_fun`	a `function` with arguments `psi1` and `psi0` specifying the estimand. Alternative, specify "ate" or "rate_ratio" as a `character` to use one of the default estimand functions. See more details in the "Estimand" section of rctglm.

estimand_fun_deriv0
: a function specifying the derivative of estimand_fun wrt. psi0. As a default the algorithm will use symbolic differentiation to automatically find the derivative from estimand_fun

estimand_fun_deriv1
: a function specifying the derivative of estimand_fun wrt. psi1. As a default the algorithm will use symbolic differentiation to automatically find the derivative from estimand_fun

cv_variance
: a logical determining whether to estimate the variance using cross-validation (see details of rctglm).

cv_variance_folds
: a numeric with the number of folds to use for cross validation if cv_variance is TRUE.

verbose
: numeric verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution (Defaults to 2, overwritable using option 'postcard.verbose' or environment variable 'R_POSTCARD_VERBOSE')

...
: Additional arguments passed to stats::glm()

Details

The procedure assumes the setup of a randomised clinical trial with observations grouped by a binary exposure_indicator variable, allocated randomly with probability exposure_prob. A GLM is fit and then used to predict the response of all observations in the event that the exposure_indicator is 0 and 1, respectively. Taking means of these predictions produce the *counterfactual means* psi0 and psi1, and an estimand r(psi0, psi1) is calculated using any specified estimand_fun.

The variance of the estimand is found by taking the variance of the influence function of the estimand. If cv_variance is TRUE, then the counterfactual predictions for each observation (which are used to calculate the value of the influence function) is obtained as out-of-sample (OOS) predictions using cross validation with number of folds specified by cv_variance_folds. The cross validation splits are performed using stratified sampling with exposure_indicator as the strata argument in rsample::vfold_cv.

Read more in vignette("model-fit").

Value

rctglm returns an object of class inheriting from "rctglm".

An object of class rctglm is a list containing the following components:

- estimand: A data.frame with plug-in estimate of estimand, standard error (SE) estimate and variance estimate of estimand
- estimand_funs: A list with
  - f: The estimand_fun used to obtain an estimate of the estimand from counterfactual means
  - d0: The derivative with respect to psi0
  - d1: The derivative with respect to psi1

- means_counterfactual: A data.frame with counterfactual means psi0 and psi1
- fitted.values_counterfactual: A data.frame with counterfactual mean values, obtained by transforming the linear predictors for each group by the inverse of the link function.
- glm: A glm object returned from running stats::glm within the procedure
- call: The matched call

Estimands

As noted in the description, psi0 and psi1 are the counterfactual means found by prediction using a fitted GLM in the binary groups defined by exposure_indicator.

Default estimand functions can be specified via "ate" (which uses the function function(psi1, psi0) psi1-psi0) and "rate_ratio" (which uses the function function(psi1, psi0) psi1/psi0). See more information on specifying the estimand_fun in vignette("model-fit").

As a default, the Deriv package is used to perform symbolic differentiation to find the derivatives of the estimand_fun.

See Also

See how to extract information using methods in rctglm_methods.

Use rctglm_with_prognosticscore() to include prognostic covariate adjustment.

See vignettes

Examples

```
# Generate some data to showcase example
n <- 100
exp_prob <- .5

dat_gaus <- glm_data(
  Y ~ 1+1.5*X1+2*A,
  X1 = rnorm(n),
  A = rbinom(n, 1, exp_prob),
  family = gaussian()
)

# Fit the model
ate <- rctglm(formula = Y ~ .,
              exposure_indicator = A,
              exposure_prob = exp_prob,
              data = dat_gaus,
              family = gaussian)

# Pull information on estimand
estimand(ate)

## Another example with different family and specification of estimand_fun
dat_binom <- glm_data(
  Y ~ 1+1.5*X1+2*A,
  X1 = rnorm(n),
```

```
  A = rbinom(n, 1, exp_prob),
  family = binomial()
)

rr <- rctglm(formula = Y ~ .,
             exposure_indicator = A,
             exposure_prob = exp_prob,
             data = dat_binom,
             family = binomial(),
             estimand_fun = "rate_ratio")

odds_ratio <- function(psi1, psi0) (psi1*(1-psi0))/(psi0*(1-psi1))
or <- rctglm(formula = Y ~ .,
             exposure_indicator = A,
             exposure_prob = exp_prob,
             data = dat_binom,
             family = binomial,
             estimand_fun = odds_ratio)
```

`rctglm_methods` *Methods for objects of class* `rctglm`

Description

Methods mostly to extract information from model fit and inference. See details for more information on each method.

Usage

```
estimand(object)

## S3 method for class 'rctglm'
estimand(object)

est(object)

## S3 method for class 'rctglm'
coef(object, ...)

## S3 method for class 'rctglm'
predict(object, ...)

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

Arguments

object	an object of class rctglm
...	additional arguments passed to methods
x	an object of class rctglm
digits	a numeric with the number of digits to display when printing

Details

The function estimand (or short-hand version est) can be used to extract a data.frame with an estimated value and standard error of the estimand.

A method for the generic coef() has been added for class rctglm, which extracts coefficient information from the underlying glm fit in the procedure. The same is true for the method defined for the predict() generic. The method for an rctglm class object calls predict.glm() on the glm fit contained within an rctglm object.

Value

estimand/est returns a data.frame with columns Estimate and Std. Error with the estimate and standard error of the estimand.

See coef() and predict.glm() for details on what is returned by the corresponding methods.

See Also

The generic rctglm() which produces rctglm class objects.

Examples

```
# Generate some data to showcase example
n <- 100
exposure_prob <- .5
dat_binom <- glm_data(
  Y ~ 1+1.5*X1+2*A,
  X1 = rnorm(n),
  A = rbinom(n, 1, exposure_prob),
  family = binomial()
)

# Fit the model
ate <- rctglm(formula = Y ~ .,
              exposure_indicator = A,
              exposure_prob = exposure_prob,
              data = dat_binom,
              family = binomial,
              estimand_fun = "ate")

print(ate)
estimand(ate)
coef(ate)
```

`rctglm_with_prognosticscore` *Use prognostic covariate adjustment when fitting an rctglm*

Description

The procedure uses fit_best_learner to fit a prognostic model to historical data and uses the model to produce counterfactual predictions as a prognostic score that is then adjusted for as a covariate in the rctglm procedure. See Powering RCTs for marginal effects with GLMs using prognostic score adjustment by Højbjerre-Frandsen et. al (2025) for more details.

Usage

```
rctglm_with_prognosticscore(
  formula,
  exposure_indicator,
  exposure_prob,
  data,
  family = gaussian,
  estimand_fun = "ate",
  estimand_fun_deriv0 = NULL,
  estimand_fun_deriv1 = NULL,
  cv_variance = FALSE,
  cv_variance_folds = 10,
  ...,
  data_hist,
  prog_formula = NULL,
  cv_prog_folds = 5,
  learners = default_learners(),
  verbose = options::opt("verbose")
)
```

Arguments

`formula`	an object of class "formula" (or one that can be coerced to that class): a symbolic description of the model to be fitted. The details of model specification are given under 'Details' in the glm documentation.
`exposure_indicator`	(name of) the *binary* variable in `data` that identifies randomisation groups. The variable is required to be binary currently.
`exposure_prob`	a `numeric` with the probability of being in "group 1" (rather than group 0) in groups defined by `exposure_indicator`.
`data`	an optional data frame, list or environment (or object coercible by as.data.frame to a data frame) containing the variables in the model. If not found in data, the variables are taken from environment(formula), typically the environment from which the function is called.

`family` a description of the error distribution and link function to be used in the model. For `glm` this can be a character string naming a family function, a family function or the result of a call to a family function. For `glm.fit` only the third option is supported. (See `family` for details of family functions.)

`estimand_fun` a `function` with arguments `psi1` and `psi0` specifying the estimand. Alternative, specify "ate" or "rate_ratio" as a `character` to use one of the default estimand functions. See more details in the "Estimand" section of rctglm.

`estimand_fun_deriv0` a `function` specifying the derivative of `estimand_fun` wrt. `psi0`. As a default the algorithm will use symbolic differentiation to automatically find the derivative from `estimand_fun`

`estimand_fun_deriv1` a `function` specifying the derivative of `estimand_fun` wrt. `psi1`. As a default the algorithm will use symbolic differentiation to automatically find the derivative from `estimand_fun`

`cv_variance` a `logical` determining whether to estimate the variance using cross-validation (see details of rctglm).

`cv_variance_folds` a `numeric` with the number of folds to use for cross validation if `cv_variance` is `TRUE`.

`...` Additional arguments passed to `stats::glm()`

`data_hist` a `data.frame` with historical data on which to fit a prognostic model

`prog_formula` an object of class "formula" (or one that can be coerced to that class): a symbolic description of the prognostic model to be fitted to `data_hist`. Passed in a list as the `preproc` argument in `fit_best_learner()`. As a default, the formula is created by modelling the response (assumed to have the same name as in `formula`) using all columns in `data_hist`.

`cv_prog_folds` a `numeric` with the number of cross-validation folds used when fitting and evaluating models

`learners` a `list` (preferably named) containing named lists of elements `model` and optionally `grid`. The `model` element should be a `parsnip` model specification, which is passed to workflowsets::workflow_set as the `model` argument, while the `grid` element is passed as the `grid` argument of workflowsets::option_add

`verbose` `numeric` verbosity level. Higher values means more information is printed in console. A value of 0 means nothing is printed to console during execution (Defaults to 2, overwritable using option 'postcard.verbose' or environment variable 'R_POSTCARD_VERBOSE')

Details

Prognostic covariate adjustment involves training a prognostic model (using fit_best_learner) on historical data (`data_hist`) to predict the response in that data.

Assuming that the historical data is representative of the comparator group in a "new" data set (group 0 of the binary `exposure_indicator` in `data`), we can use the prognostic model to predict the counterfactual outcome of all observations (including the ones in the comparator group for which the prediction of counterfactual outcome coincides with a prediction of actual outcome).

This prediction, which is called the prognostic score, is then used as an adjustment covariate in the GLM (the prognostic score is added to `formula` before calling rctglm with the modified formula).

See much more details in the reference in the description.

Value

`rctglm_with_prognosticscore` returns an object of class `rctglm_prog`, which inherits from rctglm.

An `rctglm_prog` object is a list with the same components as an rctglm object (see the `Value` section of rctglm for a breakdown of the structure), but with an additional list element of:

- `prognostic_info`: List with information about the fitted prognostic model on historical data. It has components:
  - `formula`: The `formula` with symbolic description of how the response is modelled as function of covariates in the models
  - `model_fit`: A trained `workflow` - the result of fit_best_learner
  - `learners`: A `list` of learners used for the discrete super learner
  - `cv_folds`: The amount of folds used for cross validation
  - `data`: The historical data used for cross validation when fitting and testing models

See Also

Method to extract information of the prognostic model in prog. Function used to fit the prognostic model is `fit_best_learner()`.

See `rctglm()` for the function and class this inherits from.

Examples

```
# Generate some data
n <- 100
b0 <- 1
b1 <- 1.5
b2 <- 2
W1 <- runif(n, min = 1, max = 10)
exp_prob <- .5

dat_treat <- glm_data(
  Y ~ b0+b1*log(W1)+b2*A,
  W1 = W1,
  A = rbinom (n, 1, exp_prob)
)

dat_notreat <- glm_data(
  Y ~ b0+b1*log(W1),
  W1 = W1
)

learners <- list(
  mars = list(
```

```
    model = parsnip::set_engine(
      parsnip::mars(
        mode = "regression", prod_degree = 3
      ),
      "earth"
    )
 ),
    lm = list(
      model = parsnip::set_engine(
        parsnip::linear_reg(),
        "lm"
      )
    )
)
ate <- rctglm_with_prognosticscore(
  formula = Y ~ .,
  exposure_indicator = A,
  exposure_prob = exp_prob,
  data = dat_treat,
  family = gaussian(),
  estimand_fun = "ate",
  data_hist = dat_notreat,
  learners = learners)

# Pull information on estimand
estimand(ate)
```

`repeat_power_linear` *Create data and plot power curves calculated using functions in* `power_linear()` *for a list of formulas/models*

Description

Estimate a variance for power approximation using `variance_ancova()` for each formula in `formula_list` on `train_data`. Then calculate power using the function with name specified in `power_fun` across a number of sample sizes `ns` for an assumed average treatment effect of `ate`.

Usage

```
repeat_power_linear(
  ate,
  formula_list,
  train_data,
  power_fun = c("power_gs", "power_nc"),
  ns = 10:400,
  desired_power = 0.9,
  ...
)
```

```
## S3 method for class 'postcard_rpl'
plot(x, cols = NULL, ...)
```

Arguments

ate	Passed to power_gs() or power_nc()
formula_list	a named list of formulas that are element wise passed to variance_ancova()
train_data	Passed as the data argument in variance_ancova()
power_fun	a character string with value "power_gs" or "power_nc", specifying what function in the power_linear() topic to use
ns	a numeric vector of sample sizes
desired_power	a numeric between 0 and 1 indicating the desired power level
...	Arguments passed to variance_ancova() and power_gs() or power_nc()
x	an object of class postcard_rpl created by repeat_power_linear()
cols	a (potentially named) character vector of colors for the different models in formula_list

Value

repeat_power_linear returns an object of class postcard_rpl, which is just a data.frame with a plot method defined. The plot method returns a ggplot2 object.

See Also

repeat_power_marginaleffect() for a similar implementation to iterate the process of approximating power with the power_marginaleffect()

Examples

```
train_data <- glm_data(
  Y ~ 1+1.5*log(W)+2*X,
  W = runif(1e3, min = 1, max = 10),
  X = rnorm(1e3, sd = 3)
)
rpl <- repeat_power_linear(
  ate = 0.5,
  formula_list = list("ANCOVA 1 covariate" = Y ~ X, "ANCOVA 2 covariates" = Y ~ W + X),
  train_data = train_data)

rpl_nc <- repeat_power_linear(
  ate = 0.5,
  formula_list = list("ANCOVA 1 covariate" = Y ~ X, "ANCOVA 2 covariates" = Y ~ W + X),
  train_data = train_data,
  power_fun = "power_nc",
  df = 1e3-3,
  deflation = 0.95,
  margin = -0.2,
```

```
  r = 2)

## Not run:
plot(rpl)

plot(rpl_nc)

## End(Not run)
```

`repeat_power_marginaleffect` *Create data and plot power curves calculated using* `power_marginaleffect()` *for a list of models*

Description

Iterate a process of simulating test data from `test_data_fun`, making predictions using models in `model_list`, and calculating power using `power_marginaleffect()` across a number of sample sizes `ns` and iterations `n_iter`. The results are averaged and used to create a plot of the resulting power curves.

Usage

```
repeat_power_marginaleffect(
  target_effect,
  exposure_prob,
  model_list = default_power_model_list(),
  test_data_fun = function(n) {
     glm_data(Y ~ 1 + 3 * log(W), W = stats::runif(n, min
    = 1, max = 50))
 },
  ns = seq(10, 300, 10),
  desired_power = 0.9,
  n_iter = 1,
  ...
)

## S3 method for class 'postcard_rpm'
plot(x, cols = NULL, ...)
```

Arguments

`target_effect`	Passed to `power_marginaleffect()`
`exposure_prob`	Passed to `power_marginaleffect()`
`model_list`	a named `list` of models used to get predictions on generated test data sets that are then passed to `power_marginaleffect()` as `predictions`. The elements of `model_list` need to have an existing `predict()` method. The default is an

	ANCOVA and a prognostic model fitted with `fit_best_learner()` to a simple data set of 1000 observations generated with a non-linear effect of a single covariate using `glm_data()`.
`test_data_fun`	a `function` with a single argument `n` that generates test data sets for the sample sizes `ns` specified. The default generates data using `glm_data()` with the same data generating process as the training data used to fit the default models in `model_list`.
`ns`	a `numeric` vector of sample sizes
`desired_power`	a `numeric` between 0 and 1 indicating the desired power level
`n_iter`	a `numeric` indicating a number of iterations to process and average over
`...`	additional arguments passed to `power_marginaleffect()`
`x`	an object of class `postcard_rpm` created by `repeat_power_marginaleffect()`
`cols`	a (potentially named) `character` vector of colors for the different models in `model_list`

Value

`repeat_power_marginal` returns an object of class `postcard_rpm`, which is just a `data.frame` with a `plot` method defined. The `plot` method returns a `ggplot2` object.

See Also

`repeat_power_linear()` for a similar implementation to iterate the process of approximating power with the functions in `power_linear()`

Examples

```
# Note everything is wrapped in dontrun to avoid long runtimes of examples (tests are
# still in place). Reduce the number of sample sizes and/or iterations to avoid long
# runtimes
## Not run:
# A simple use case with default models and test data (we run only with a few sample
# sizes to reduce runtime of examples)
rpm <- repeat_power_marginaleffect(
  target_effect = 0.9,
  exposure_prob = 0.5
)
plot(rpm)

################################
# Create model from a poisson family and estimate the power of rate ratio with
# several arguments passed to power_marginaleffect
################################
b1 <- 0.9
b2 <- 0.2
b3 <- -0.4
b4 <- -0.6

train_pois <- glm_data(
```

```
  Y ~ b1*log(X1)+b2*X2+b3*X3+b4*X2*X3,
  X1 = runif(1e3, min = 1, max = 10),
  X2 = rnorm(1e3),
  X3 = rgamma(1e3, shape = 1),
  family = poisson()
)

# Define models to compare fit to training data
ancova_prog_list <- list(
ANCOVA = glm(Y ~ X1 + X2 + X3, data = train_pois, family = poisson),
"ANCOVA with prognostic score" = fit_best_learner(list(mod = Y ~ X1 + X2 + X3), data = train_pois)
)

# Create a function that produces data to predict on
test_pois_fun <- function(n) {
 glm_data(
   Y ~ b1*log(X1)+b2*X2+b3*X3+b4*X2*X3,
   X1 = runif(n, min = 1, max = 10),
   X2 = rnorm(n),
   X3 = rgamma(n, shape = 1),
   family = poisson()
 )
}

# Specify a bunch of different arguments that are passed to power_marginaleffect()
## Run for 2 sample sizes to reduce runtime
rpm_rr <- repeat_power_marginaleffect(
  model_list = ancova_prog_list,
  test_data_fun = test_pois_fun,
  ns = seq(100, 200), n_iter = 1,
  var1 = function(var0) 1.1 * var0,
  kappa1_squared = function(kap0) 1.1 * kap0,
  estimand_fun = "rate_ratio",
  target_effect = 1.4,
  exposure_prob = 1/2,
  margin = 0.8
)
plot(rpm_rr2)

## End(Not run)
```

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