

# Package ‘propertee’

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**Version** 1.0.4

**Title** Standardization-Based Effect Estimation with Optional Prior Covariance Adjustment

**Description** The Prognostic Regression Offsets with Propagation of EErrors (for Treatment Effect Estimation) package facilitates direct adjustment for experiments and observational studies that is compatible with a range of study designs and covariance adjustment strategies. It uses explicit specification of clusters, blocks and treatment allocations to furnish probability of assignment-based weights targeting any of several average treatment effect parameters, and for standard error calculations reflecting these design parameters. For covariance adjustment of its Hajek and (one-way) fixed effects estimates, it enables offsetting the outcome against predictions from a dedicated covariance model, with standard error calculations propagating error as appropriate from the covariance model.

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 'SandwichLayer.R' 'block\_center\_residuals.R'  
 'SandwichLayerVariance.R' 'StudySpecificationConverters.R'  
 'StudySpecificationStructure.R' 'StudySpecificationUtilities.R'  
 'WeightedStudySpecification.R' 'areg.center.R' 'confint\_lm.R'  
 'teeMod.R' 'as.lmitt.R' 'as\_data\_frame.R' 'assigned.R'  
 'bread.R' 'c\_weightedStudySpecification.R' 'cov\_adj.R' 'data.R'  
 'dichotomy.R' 'expand.model.frame\_tee.R' 'get\_cov\_adj.R'  
 'get\_data\_from\_model.R' 'get\_spec.R' 'glmrobMethods.R'  
 'lmitt.R' 'lmrob\_methods.R' 'merge\_preserve\_order.R'  
 'propertee\_package.R' 'specification\_table.R'  
 'summary.StudySpecification.R' 'summary.teeMod.R' 'update\_by.R'  
 'validWeights.R' 'weights\_internal.R' 'weights\_exported.R'

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+, WeightedStudySpecification, numeric-method

*WeightedStudySpecification Operations***Description**

Algebraic operators on WeightedStudySpecification objects and numeric vectors. WeightedStudySpecifications do not support addition or subtraction.

**Usage**

```
## S4 method for signature 'WeightedStudySpecification,numeric'
e1 + e2
```

```
## S4 method for signature 'numeric,WeightedStudySpecification'
e1 + e2
```

```
## S4 method for signature 'WeightedStudySpecification,numeric'
e1 - e2
```

```
## S4 method for signature 'numeric,WeightedStudySpecification'
e1 - e2
```

```
## S4 method for signature 'WeightedStudySpecification,numeric'
e1 * e2
```

```
## S4 method for signature 'numeric,WeightedStudySpecification'
e1 * e2

## S4 method for signature 'WeightedStudySpecification,numeric'
e1 / e2

## S4 method for signature 'numeric,WeightedStudySpecification'
e1 / e2
```

### Arguments

e1, e2                      WeightedStudySpecification or numeric objects

### Details

These are primarily used to either combine weights via multiplication, or to invert weights. Addition and subtraction are not supported and will produce errors.

### Value

a WeightedStudySpecification object

---

as.lmitt	<i>Convert lm object into teeMod</i>
----------	--------------------------------------

---

### Description

Converts the output of `lm()` into a `teeMod` object, for standard errors that account for block and cluster information carried with the `lm`'s weights, and/or an offset incorporating predictions of the outcome from a separate model.

### Usage

```
as.lmitt(x, specification = NULL)
```

```
as.teeMod(x, specification = NULL)
```

### Arguments

x	lm object with weights containing a WeightedStudySpecification, or an offset from <code>cov_adj()</code> .
specification	Optional, explicitly specify the StudySpecification to be used. If the StudySpecification is specified elsewhere in x (e.g. passed as an argument to any of <code>ate()</code> , <code>ett()</code> , <code>cov_adj()</code> or <code>assigned()</code> ) it will be found automatically and does not need to be passed here as well. If different StudySpecification objects are passed (either through the <code>lm</code> in weights or covariance adjustment, or through this argument), an error will be produced.

## Details

The formula with which `x` was created must include a treatment identifier (e.g. `assigned()`). If a model-based offset is incorporated, the model's predictions would have to have been extracted using `cov_adj()` as opposed to `predict()` in order for `teeMod` standard error calculations to reflect propagation of error from these predictions. This mechanism only supports treatment main effects: to estimate interactions of treatment assignment with a moderator variable, use `lmitt()` instead of `lm()` followed by `as.lmitt()`.

## Value

`teeMod` object

---

<code>as.SandwichLayer</code>	<i>Convert a <code>PreSandwichLayer</code> to a <code>SandwichLayer</code> with a <code>StudySpecification</code> object</i>
-------------------------------	--

---

## Description

`as.SandwichLayer()` uses the `StudySpecification` object passed to the `specification` argument to populate the slots in a `SandwichLayer` object that a `PreSandwichLayer` does not have sufficient information for.

## Usage

```
as.SandwichLayer(x, specification, by = NULL, Q_data = NULL)
```

## Arguments

<code>x</code>	a <code>PreSandwichLayer</code> object
<code>specification</code>	a <code>StudySpecification</code> object
<code>by</code>	optional; a string or named vector of unique identifier columns in the data used to create <code>specification</code> and the data used to fit the covariance adjustment model. Default is <code>NULL</code> , in which case unit of assignment columns are used for identification (even if they do not uniquely identify units of observation). If a named vector is provided, names should represent variables in the data used to create <code>specification</code> , while values should represent variables in the covariance adjustment data.
<code>Q_data</code>	dataframe of direct adjustment sample, which is needed to generate the keys slot of the <code>SandwichLayer</code> object. Defaults to <code>NULL</code> , in which case if <code>by</code> is <code>NULL</code> , the data used to create <code>specification</code> is used, and if <code>by</code> is not <code>NULL</code> , appropriate data further up the call stack (passed as arguments to <code>cov_adj()</code> or <code>lmitt.formula()</code> , for example) is used.

## Value

a `SandwichLayer` object

assigned

*Obtain Treatment from StudySpecification***Description**

When passing a `lm` object to `lmitt()`, extract and use the treatment variable specified in the `StudySpecification`.

**Usage**

```
assigned(specification = NULL, data = NULL, dichotomy = NULL)
```

```
adopters(specification = NULL, data = NULL, dichotomy = NULL)
```

```
a.(specification = NULL, data = NULL, dichotomy = NULL)
```

```
z.(specification = NULL, data = NULL, dichotomy = NULL)
```

**Arguments**

<code>specification</code>	Optional <code>StudySpecification</code> . If the <code>StudySpecification</code> can't be identified in the model (usually because neither <code>weights(ate())</code> or <code>ett()</code> nor a covariate adjustment model ( <code>cov_adj()</code> ) are found), the <code>StudySpecification</code> can be passed directly.
<code>data</code>	Optional data set. By default <code>assigned()</code> will attempt to identify the appropriate data, if this fails (or you want to overwrite it), you can pass the data here.
<code>dichotomy</code>	optional; a formula defining the dichotomy of the treatment variable if it isn't already 0/1. See details for more information. If <code>ett()</code> or <code>ate()</code> is called within a <code>lmitt()</code> call that specifies a dichotomy argument, that dichotomy will be used if the argument here has not been specified.

**Details**

When passing a `lm` object to `lmitt()`, the treatment variable in the formula passed to `lm()` needs to be identifiable. Rather than placing the treatment variable directly in the formula, use one of these functions, to allow `lmitt()` to identify the treatment variable.

To keep the formula in the `lm()` call concise, instead of passing `specification` and `data` arguments to these functions, one can pass a `WeightedStudySpecification` object to the `weights` argument of the `lm()` call or a `SandwichLayer` object to the `offset` argument.

Alternatively, you can pass the `specification` and `data` arguments.

While `assigned()` can be used in any situation, it is most useful for scenarios where the treatment variable is non-binary and the `StudySpecification` contains a `Dichotomy`. For example, say `q` is a 3-level ordinal treatment variable, and the binary comparison of interest is captured in `dichotomy = q == 3 ~ q < 3`. If you were to fit a model including `q` as a predictor, e.g. `lm(y ~ q, ...)`, `lm` would treat `q` as the full ordinal variable. On the other hand, by calling `lm(y ~ assigned(), weights =`

`ate(spec), ...)`, `assigned()` will generate the appropriate binary variable to allow estimation of treatment effects.

If called outside of a model call and without a `data` argument, this will extract the treatment from the specification. If this is the goal, the `treatment()` function is better suited for this purpose.

### Value

The treatment variable to be placed in the regression formula.

### Examples

```
data(simdata)
spec <- obs_spec(z ~ uoa(uoa1, uoa2), data = simdata)
mod <- lm(y ~ assigned(), data = simdata, weights = ate(spec))
lmittmod <- lmitt(mod)
summary(lmittmod, vcov.type = "CR0")
```

---

as_rct_spec	<i>Convert StudySpecification between types</i>
-------------	---

---

### Description

Convert a `StudySpecification` between a observational study, a randomized control trial, and a regression discontinuity (created from `obs_spec`, `rct_spec` and `rd_spec` respectively).

### Usage

```
as_rct_spec(StudySpecification, ..., loseforcing = FALSE)
```

```
as_obs_spec(StudySpecification, ..., loseforcing = FALSE)
```

```
as_rd_spec(StudySpecification, data, ..., forcing)
```

### Arguments

<code>StudySpecification</code>	a <code>StudySpecification</code> to convert
<code>...</code>	Ignored.
<code>loseforcing</code>	converting from RD to another <code>StudySpecification</code> type will error to avoid losing the forcing variable. Setting <code>loseforcing = TRUE</code> allows the conversion to automatically drop the forcing variable. Default <code>FALSE</code> .
<code>data</code>	converting to an RD requires adding a forcing variable, which requires access to the original data.
<code>forcing</code>	converting to an RD requires adding a forcing variable. This should be entered as a formula which would be passed to <code>update</code> , e.g. <code>forcing = . ~ . + forcing(forcevar)</code> .

**Value**

StudySpecification of the updated type

**Examples**

```
spec <- rct_spec(z ~ unit_of_assignment(uoa1, uoa2), data = simdata)
spec
as_obs_spec(spec)
as_rd_spec(spec, simdata, forcing = ~ . + forcing(force))
spec2 <- rd_spec(o ~ uoa(uoa1, uoa2) + forcing(force), data = simdata)
spec2
# as_rct_spec(spec2) # this will produce an error
as_rct_spec(spec2, loseforcing = TRUE)
```

---

bread.teeMod

---

*Extract bread matrix from a teeMod model fit*


---

**Description**

An S3method for `sandwich::bread` that extracts the bread of the direct adjustment model sandwich covariance matrix.

**Usage**

```
## S3 method for class 'teeMod'
bread(x, ...)
```

**Arguments**

x	a fitted teeMod model
...	arguments passed to methods

**Details**

This function is a thin wrapper around `.get_tilde_a22_inverse()`.

**Value**

A variance-covariance matrix with row and column entries for the estimated coefficients in x, the marginal mean outcome in the control condition, the marginal mean offset in the control condition (if an offset is provided), and if a moderator variable is specified in the formula for x, the mean interaction in the control condition of the outcome and offset with the moderator variable



---

c, WeightedStudySpecification-method  
*Concatenate weights*

---

## Description

Given several variations of weights generated from a single StudySpecification, combine into a single weight.

## Usage

```
## S4 method for signature 'WeightedStudySpecification'
c(x, ..., warn_dichotomy_not_equal = FALSE)
```

## Arguments

**x** .. a WeightedStudySpecification object, typically created from `ate()` or `ett()`

**...** any number of additional WeightedStudySpecification objects with equivalent StudySpecification to x and each other

**warn\_dichotomy\_not\_equal** if FALSE (default), WeightedStudySpecifications are considered equivalent even if their dichotomy differs. If TRUE, a warning is produced.

## Details

Concatenating WeightedStudySpecification objects with `c()` requires both individual WeightedStudySpecification objects to come from the same StudySpecification and have the same target (e.g all created with `ate()` or all created with `ett()`, no mixing-and-matching). All arguments to `c()` must be WeightedStudySpecification.

WeightedStudySpecification objects may be concatenated together even without having the same @dichotomy slot. This procedure only prompts a warning for differing dichotomies if the argument `warn_dichotomy_not_equal` is set to TRUE.

## Value

A numeric vector with the weights concatenated in the input order.

## Examples

```
data(simdata)
spec <- rct_spec(z ~ unit_of_assignment(uoa1, uoa2), data = simdata)
w1 <- ate(spec, data = simdata[1:30,])
w2 <- ate(spec, data = simdata[31:40,])
w3 <- ate(spec, data = simdata[41:50,])
c_w <- c(w1, w2, w3)
c(length(w1), length(w2), length(w3), length(c_w))
```

```
spec <- rct_spec(dose ~ unit_of_assignment(uoa1, uoa2), data = simdata)
w1 <- ate(spec, data = simdata[1:10, ], dichotomy = dose >= 300 ~ .)
w2 <- ate(spec, data = simdata[11:30, ], dichotomy = dose >= 200 ~ .)
w3 <- ate(spec, data = simdata[31:50, ], dichotomy = dose >= 100 ~ .)
c_w <- c(w1, w2, w3)
```

---

confint.teeMod	<i>Confidence intervals with standard errors provided by</i> vcov.teeMod()
----------------	---

---

## Description

An S3method for `stats::confint` that uses standard errors computed using `vcov.teeMod()`. Additional arguments passed to this function, such as `cluster` and `type`, specify the arguments of the `vcov.teeMod()` call.

## Usage

```
## S3 method for class 'teeMod'
confint(object, parm, level = 0.95, ...)
```

## Arguments

<code>object</code>	a fitted <code>teeMod</code> model
<code>parm</code>	a specification of which parameters are to be given confidence intervals, either a vector of numbers or a vector of names. If missing, all parameters are considered.
<code>level</code>	the confidence level required.
<code>...</code>	additional arguments to pass to <code>vcov.teeMod()</code>

## Details

Rather than call `stats::confint.lm()`, `confint.teeMod()` calls `.confint_lm()`, a function internal to the `proptee` package that ensures additional arguments in the `...` of the `confint.teeMod()` call are passed to the internal `vcov()` call.

## Value

A matrix (or vector) with columns giving lower and upper confidence limits for each parameter. These will be labelled as  $(1-\text{level})/2$  and  $1 - (1-\text{level})/2$  in % (by default 2.5% and 97.5%)

---

cov_adj	<i>Covariance adjustment of teeMod model estimates</i>
---------	--

---

### Description

`cov_adj()` takes a fitted covariance model and returns the information necessary for adjusting direct adjustment model estimates and associated standard errors for covariates. Standard errors will reflect adjustments made to the outcomes as well as contributions to sampling variability arising from the estimates of the covariance adjustment model coefficients.

### Usage

```
cov_adj(model, newdata = NULL, specification = NULL, by = NULL)
```

### Arguments

<code>model</code>	any model that inherits from a <code>glm</code> , <code>lm</code> , or <code>robustbase::lmrob</code> object
<code>newdata</code>	a dataframe of new data. Default is <code>NULL</code> , in which case a dataframe is sought from higher up the call stack.
<code>specification</code>	a <code>StudySpecification</code> object. Default is <code>NULL</code> , in which case a <code>StudySpecification</code> object is sought from higher up the call stack.
<code>by</code>	optional; a string or named vector of unique identifier columns in the data used to create <code>specification</code> and the data used to fit the covariance adjustment model. Default is <code>NULL</code> , in which case unit of assignment columns are used for identification (even if they do not uniquely identify units of observation). If a named vector is provided, names should represent variables in the data used to create <code>specification</code> , while values should represent variables in the covariance adjustment data.

### Details

Prior to generating adjustments, `cov_adj()` identifies the treatment variable specified in the `StudySpecification` object passed to `specification` and replaces all values with a reference level. If the treatment has logical type, this reference level is `FALSE`, and if it has numeric type, this is the smallest non-negative value (which means 0 for 0/1 binary). Factor treatments are not currently supported for `StudySpecification` objects.

The values of the output vector represent adjustments for the outcomes in `newdata` if `newdata` is provided; adjustments for the outcomes in the data used to fit a `teeMod` model if `cov_adj()` is called within the `offset` argument of the model fit; or they are the fitted values from `model` if no relevant dataframe can be extracted from the call stack. The length of the output of `cov_adj()` will match the number of rows of the dataframe used.

### Value

A `SandwichLayer` if `specification` is not `NULL` or a `StudySpecification` object is found in the call stack, otherwise a `PreSandwichLayer` object

## Examples

```
data("STARplus")

##' A prognostic model fitted to experimental + non-experimental controls
y0hat_read <- lm(read_yr1 ~ gender*dob +dobNA + race,
                 data = STARplus,
                 subset = cond_at_entry!="small")

STARspec <- rct_spec(cond_at_entry ~ unit_of_assignment(stdntid) +
                    block(grade_at_entry, school_at_entry),
                    subset=!is.na(grade_at_entry),# excludes non-experimentals
                    data = STARplus)
ett_wts <- ett(STARspec, data = STARplus,
               dichotomy= cond_at_entry=="small" ~.)

ett_read <- lm(read_yr1 ~ assigned(dichotomy= cond_at_entry=="small" ~.),
               offset = cov_adj(y0hat_read),
               data = STARplus,
               weights = ett_wts)
coef(ett_read)
ett_read |> as.lmitt() # brings in control-group means of outcome, predictions

ate_read <- lmitt(read_yr1 ~ 1, STARspec, STARplus,
                  dichotomy= cond_at_entry=="small" ~.,
                  offset = cov_adj(y0hat_read),
                  weights = "ate")
show(ate_read)
vcov(ate_read, type = "HC0", cov_adj_rcorrect = "HC0") |> unname()

ate_read_loc <-
  lmitt(read_yr1 ~ race, STARspec, STARplus,
        dichotomy= cond_at_entry=="small" ~.,
        offset = cov_adj(y0hat_read, newdata = STARplus),
        weights = "ate")
show(ate_read_loc)
```

---

estfun.glmrob

---

*Extract empirical estimating equations from a glmrob model fit*


---

## Description

Extract empirical estimating equations from a glmrob model fit

Extract bread matrix from an lmrob() fit

## Usage

```
## S3 method for class 'glmrob'
estfun(x, ...)
```

```
## S3 method for class 'glmrob'
bread(x, ...)
```

### Arguments

x                      a fitted lmrob object  
 ...                    arguments passed to methods

### Value

matrix, estimating functions evaluated at data points and fitted parameters  
 matrix, inverse Hessian of loss as evaluated at fitted parameters

---

estfun.lmrob	<i>Generate matrix of estimating equations for lmrob() fit</i>
--------------	--

---

### Description

Generate matrix of estimating equations for lmrob() fit  
 Extract bread matrix from an lmrob() fit

### Usage

```
## S3 method for class 'lmrob'
estfun(x, ...)

## S3 method for class 'lmrob'
bread(x, ...)
```

### Arguments

x                      An lmrob object produced using an MM/SM estimator chain  
 ...                    Additional arguments to be passed to bread

### Details

This is part of a workaround for an issue in the robustbase code affecting sandwich covariance estimation. The issue in question is issue #6471, robustbase project on R-Forge. This function contributes to providing sandwich estimates of covariance-adjusted standard errors for robust linear covariance adjustment models.

This is part of a workaround for an issue in the robustbase code affecting sandwich covariance estimation. The issue in question is issue #6471, robustbase project on R-Forge. This function contributes to providing sandwich estimates of covariance-adjusted standard errors for robust linear covariance adjustment models.

**Value**

A  $n \times (p+1)$  matrix where the first column corresponds to the scale estimate and the remaining  $p$  columns correspond to the coefficients

A  $p \times (p+1)$  matrix where the first column corresponds to the scale estimate and the remaining  $p$  columns correspond to the coefficients

**Author(s)**

Ben B. Hansen

---

estfun.teeMod

*Extract empirical estimating equations from a teeMod model fit*

---

**Description**

An S3method for sandwich::estfun for producing a matrix of contributions to the direct adjustment estimating equations.

**Usage**

```
## S3 method for class 'teeMod'
estfun(x, ...)
```

**Arguments**

x	a fitted teeMod model
...	arguments passed to methods, most importantly those that define the bias corrections for the residuals of x and, if applicable, a fitted_covariance_model stored in its offset

**Details**

If a prior covariance adjustment model has been passed to the offset argument of the teeMod model using cov\_adj(), estfun.teeMod() incorporates contributions to the estimating equations of the covariance adjustment model.

The covariance adjustment sample may not fully overlap with the direct adjustment sample, in which case estfun.teeMod() returns a matrix with the same number of rows as the number of unique units of observation used to fit the two models. Uniqueness is determined by matching units of assignment used to fit the covariance adjustment model to units of assignment in the teeMod model's StudySpecification slot; units of observation within units of assignment that do not match are additional units that add to the row count.

Theby argument in cov\_adj() can provide a column or a pair of columns (a named vector where the name specifies a column in the direct adjustment sample and the value a column in the covariance adjustment sample) that uniquely specifies units of observation in each sample. This information

can be used to align each unit of observation's contributions to the two sets of estimating equations. If no `by` argument is provided and units of observation cannot be uniquely specified, contributions are aligned up to the unit of assignment level. If standard errors are clustered no finer than that, they will provide the same result as if each unit of observation's contributions were aligned exactly.

This method incorporates bias corrections made to the residuals of `x` and, if applicable, the covariance model stored in its `offset`. When its crossproduct is taken (perhaps after suitable summing across rows within clusters), it provides a heteroskedasticity- (or cluster-) robust estimate of the meat matrix of the variance-covariance of the parameter estimates in `x`.

### Value

An  $n \times k$  matrix of empirical estimating equations for `x`. `k` includes the model intercept, main effects of treatment and moderator variables, any moderator effects, and marginal and conditional means of the outcome (and `offset`, if provided) in the control condition. See Details for definition of `n`.

---

 ett

---

*Generate Direct Adjusted Weights for Treatment Effect Estimation*


---

### Description

These should primarily be used inside models. See Details.

### Usage

```
ett(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
att(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
ate(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
etc(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
atc(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
ato(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
olw(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
owt(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
pwt(specification = NULL, dichotomy = NULL, by = NULL, data = NULL)
```

## Arguments

specification	optional; a <code>StudySpecification</code> object created by one of <code>rct_spec()</code> , <code>rd_spec()</code> , or <code>obs_spec()</code> .
dichotomy	optional; a formula defining the dichotomy of the treatment variable if it isn't already 0/1. See details for more information. If <code>ett()</code> or <code>ate()</code> is called within a <code>lmitt()</code> call that specifies a dichotomy argument, that dichotomy will be used if the argument here has not been specified.
by	optional; named vector or list connecting names of unit of assignment/ variables in specification to unit of assignment/unitid/cluster variables in data. Names represent variables in the <code>StudySpecification</code> ; values represent variables in the data. Only needed if variable names differ.
data	optional; the data for the analysis to be performed on. May be excluded if these functions are included as the <code>weights</code> argument of a model.

## Details

These functions should primarily be used in the `weight` argument of `lmitt()` or `lm()`. All arguments are optional if used within those functions. If used on their own, `specification` and `data` must be provided.

- `ate` - Average treatment effect. Aliases: `ate()`.
- `ett` - Effect of treatment on the treated. Aliases: `ett()`, `att()`.
- `etc` - Effect of treatment on controls. Aliases: `etc()`, `atc()`.
- `ato` - Overlap-weighted average effect. Aliases: `ato()`, `olw`, `owt`, `pwt`.

In a `StudySpecification` with blocks, the weights are generated as a function of the ratio of the number of treated units in a block versus the total number of units in a block.

In any blocks where that ratio is 0 or 1 (that is, all units in the block have the same treatment status), the weights will be 0. In effect this removes from the target population any block in which there is no basis for estimating either means under treatment or means under control.

If `block` is missing for a given observation, a weight of 0 is applied.

A dichotomy is specified by a formula consisting of a conditional statement on both the left-hand side (identifying treatment levels associated with "treatment") and the right hand side (identifying treatment levels associated with "control"). For example, if your treatment variable was called `dose` and doses above 250 are considered treatment, you might write:

```
ate(..., dichotomy = dose > 250 ~ dose <= 250
```

The period (.) can be used to assign all other units of assignment. For example, we could have written the same treatment regime as either

```
etc(..., dichotomy = dose > 250 ~ .
```

or

```
olw(..., dichotomy = . ~ dose <= 250
```

The dichotomy formula supports Relational Operators (see [Comparison](#)), Logical Operators (see [Logic](#)), and `%in%` (see [match\(\)](#)).



The conditionals need not assign all values of treatment to control or treatment, for example, `dose > 300 ~ dose < 200` does not assign `200 <= dose <= 300` to either treatment or control. This would be equivalent to manually generating a binary variable with NA whenever dose is between 200 and 300. Standard errors will reflect the sizes of the comparison groups specified by the dichotomy.

Tim Lyncurgus contributed code for the computation of weights. The ‘overlap weight’ concept is due to Li, Morgan and Zaslavsky (2018), although the current implementation differs from that discussed in their paper in that it avoids estimated propensity scores.

## Value

a `WeightedStudySpecification` object, which is a vector of numeric weights

## References

Li, Fan, Kari Lock Morgan, and Alan M. Zaslavsky. "Balancing covariates via propensity score weighting." *Journal of the American Statistical Association* 113, no. 521 (2018): 390-400.

## Examples

```
data(simdata)
spec <- rct_spec(z ~ unit_of_assignment(uoa1, uoa2), data = simdata)
summary(lmitt(y ~ 1, data = simdata, specification = spec, weights = ate()), vcov.type = "CR0")
```

---

get\_structure

StudySpecification *Structure Information*

---

## Description

Obtaining a `data.frame` which encodes the specification information.

## Usage

```
get_structure(specification)

## S4 method for signature 'StudySpecificationStructure'
show(object)
```

## Arguments

`specification` a `StudySpecification` object

`object` a `StudySpecificationStructure` object, typically the output of `get_structure`

## Value

A `StudySpecificationStructure` object containing the structure of the specification as a `data.frame`.

## Examples

```
data(simdata)
spec <- rct_spec(z ~ uoa(uoa1, uoa2) + block(bid), data = simdata)
get_structure(spec)
```

---

GV_data	<i>Cluster-randomized experiment data on voter turnout in cable system markets</i>
---------	--

---

## Description

This dataset is a toy example derived from a cluster-randomized field experiment that evaluates the effect of “Rock the Vote” TV advertisements on voter turnout rate. The original study included 23,869 first-time voters across 85 cable television markets in 12 states. These markets were grouped into matched sets based on their past voter turnout rates and then randomly assigned to either a treatment or control condition. This toy dataset is constructed by randomly sampling 10% of individuals from selected cable television markets in the original dataset.

## Usage

```
GV_data
```

## Format

A data.frame with 248 rows and 7 columns.

- age Age of participant
- vote\_04 Outcome variable indicating whether participant voted
- tv\_company Cable system serving participant’s residential area
- treatment Binary variable denoting treatment assignment
- pairs A numeric indicator for the strata or matched pair group to which a cable system belongs (1-3)
- population\_size Total population size of residential area served by cable system
- sample\_size Number of individuals sampled from the cable system cluster

## Details

The original dataset was drawn from a randomized controlled trial in which 85 cable system areas were first grouped into 40 matched sets based on historical voter turnout. Within each matched set, one cable system area was randomly assigned to the treatment condition, while the others served as controls.

This toy dataset includes a subset of the original replication data, specifically individuals from matched sets 1–3, which encompass 7 of the 85 cable system areas. Within these selected clusters, a 10% random sample of individuals was taken.

The fuller Green-Vavreck dataset that this derives from bears a Creative Commons BY-NC-ND license (v3.0) and is housed in Yale University’s Institution for Social and Policy Studies (ID: D005).

**Source**

<https://isps.yale.edu/research/data/d005>

**References**

Green, Donald P. & Lynn Vavreck (2008) "Analysis of Cluster-Randomized Experiments: A Comparison of Alternative Estimation Approaches." *Political Analysis* 16(2):138-152.

---

has_binary_treatment	<i>Check whether treatment stored in a StudySpecification object is binary</i>
----------------------	--

---

**Description**

Check whether treatment stored in a StudySpecification object is binary

**Usage**

```
has_binary_treatment(spec)
```

**Arguments**

spec	StudySpecification object
------	---------------------------

**Value**

logical vector of length 1

---

identical_StudySpecifications	<i>Test equality of two StudySpecification objects</i>
-------------------------------	--

---

**Description**

Check whether two StudySpecification objects are identical.

**Usage**

```
identical_StudySpecifications(x, y)
```

**Arguments**

x	A StudySpecification object.
y	A StudySpecification object.

**Value**

Logical, are x and y identical?

---

`identify_small_blocks` *Identify fine strata*

---

### Description

Identify blocks in a `StudySpecification` with exactly one treated or one control unit of assignment.

### Usage

```
identify_small_blocks(spec)
```

### Arguments

`spec`                      A `StudySpecification` object.

### Value

Logical vector with length given by the number of blocks in `StudySpecification`

---

`lmitt`                      *Linear Model for Intention To Treat*

---

### Description

Generates a linear model object to estimate a treatment effect, with proper estimation of variances accounting for the study specification.

### Usage

```
lmitt(obj, specification, data, ...)
```

```
## S3 method for class 'formula'
```

```
lmitt(
  obj,
  specification,
  data,
  absorb = FALSE,
  offset = NULL,
  weights = NULL,
  ...
)
```

```
## S3 method for class 'lm'
```

```
lmitt(obj, specification = NULL, ...)
```

## Arguments

<code>obj</code>	A formula or a <code>lm</code> object. See Details.
<code>specification</code>	The <code>StudySpecification</code> to be used. Alternatively, a formula creating a specification. (Of the type of that would be passed as the first argument to <code>rd_spec()</code> , <code>rct_spec()</code> , or <code>obs_spec()</code> , with the difference that <code>cluster()</code> , <code>uoa()</code> and <code>unit_of_assignment()</code> terms can be omitted when each row of data represents a distinct unit of assignment.) If the formula includes a <code>forcing()</code> element, an RD specification is created. Otherwise an observational specification is created. An RCT specification must be created manually using <code>rct_spec()</code> .
<code>data</code>	A <code>data.frame</code> such as would be passed into <code>lm()</code> .
<code>...</code>	Additional arguments passed to <code>lm()</code> and other functions. An example of the latter is <code>dichotomy=</code> , a formula passed to <code>assigned()</code> and, as appropriate, <code>ate()</code> , <code>att()</code> , <code>atc()</code> or <code>ato()</code> . It is used to dichotomize a non-binary treatment variable in specification. See the Details section of the <code>ate()</code> help page for examples.
<code>absorb</code>	If <code>TRUE</code> , fixed effects are included for blocks identified in the <code>StudySpecification</code> . Excluded in <code>FALSE</code> . Default is <code>FALSE</code> . The estimates of these fixed effects are suppressed from the returned object.
<code>offset</code>	Offset of the kind which would be passed into <code>lm()</code> . Ideally, this should be the output of <code>cov_adj()</code> .
<code>weights</code>	Which weights should be generated? Options are <code>"ate"</code> or <code>"ett"</code> . Alternatively, the output of a manually run <code>ate()</code> or <code>ett()</code> can be used.

## Details

The first argument to `lmitt()` should be a formula specifying the outcome on the left hand side. The right hand side of the formula can be any of the following:

- 1: Estimates a main treatment effect.
- a subgroup variable: Estimates a treatment effect within each level of your subgrouping variable.
- a continuous moderator: Estimates a main treatment effect as well as a treatment by moderator interaction. The moderator is not automatically centered.

Alternatively, `obj` can be a pre-created `lm` object. No modification is made to the formula of the object. See the help for `as.lmitt()` for details of this conversion.

The `lmitt()` function's `subset=` argument governs the subsetting of data prior to model fitting, just as with `lm()`. Functions such as `rct_spec()` that create `StudySpecifications` also take an optional `subset=` argument, but its role differs from that of the `subset=` argument of `lm()` or `lmitt()`. The `subset=` argument when creating a `StudySpecification` restricts the data used to generate the `StudySpecification`, but has no direct impact on the future `lm()` or `lmitt()` calls using that `StudySpecification`. (It can have an indirect impact by excluding particular units from receiving a treatment assignment or weight. When treatment assignments or weights are reconstructed from the `StudySpecification`, these units will receive `NA`s, and will be excluded from the `lm()` or `lmitt()` fit under typical `na.action` settings.)

To avoid variable name collision, the treatment variable defined in the specification will have a "." appended to it. For example, if you request a main treatment effect (with a formula of  $\sim 1$ ) with a treatment variable named "txt", you can obtain its estimate from the returned teeMod object via `$coefficients["txt."]`.

`lmitt()` will produce a message if the StudySpecification designates treatment assignment by block but the blocking structure appears not to be reflected in the weights, nor in a block fixed effect adjustment (via `absorb=TRUE`). While not an error, this is at odds with intended uses of `propertee`, so `lmitt()` flags it as a potential oversight on the part of the analyst. To disable this message, run `options("propertee_message_on_unused_blocks" = FALSE)`.

`lmitt()` returns objects of class 'teeMod', for Treatment Effect Estimate Model, extending the `lm` class to add a summary of the response distribution under control (the coefficients of a controls-only regression of the response on an intercept and any moderator variable). `teeMod` objects also record the underlying StudySpecification and information about any externally fitted models `mod` that may have been used for covariance adjustment by passing `offset=cov_adj(mod)`. In the latter case, responses are offsetted by predictions from `mod` prior to treatment effect estimation, but estimates of the response variable distribution under control are calculated without reference to `mod`.

The response distribution under control is also characterized when treatment effects are estimated with block fixed effects, i.e. for `lmitt()` with a formula first argument with option `absorb=TRUE`. Here as otherwise, the supplementary coefficients describe a regression of the response on an intercept and moderator variables, to which only control observations contribute; but in this case the weights are modified for this supplementary regression. The treatment effect estimates adjusted for block fixed effects can be seen to coincide with estimates calculated without block effect but with weights multiplied by an additional factor specific to the combination of block and treatment condition. For block  $s$  containing units with weights  $w_i$  and binary treatment assignments  $z_i$ , define  $\hat{\pi}_s$  by  $\hat{\pi}_s \sum_s w_i = \sum_s z_i w_i$ . If  $\hat{\pi}_s$  is 0 or 1, the block doesn't contribute to effect estimation and the additional weighting factor is 0; if  $0 < \hat{\pi}_s < 1$ , the additional weighting factor is  $1 - \hat{\pi}_s$  for treatment group members and  $\hat{\pi}_s$  for controls. When estimating a main effect only or a main effect with continuous moderator, supplementary coefficients under option `absorb=TRUE` reflect regressions with additional weighting factor equal to 0 or  $\hat{\pi}_s$ , respectively, for treatment or control group members of block  $s$ . With a categorical moderator and `absorb=TRUE`, this additional weighting factor determining supplementary coefficients is calculated separately for each level  $\ell$  of the moderator variable, with the sums defining  $\hat{\pi}_{s\ell}$  restricted not only to block  $s$  but also to observations with moderator equal to  $\ell$ .

## Value

teeMod object (see Details)

## Examples

```
data(simdata)
spec <- rct_spec(z ~ cluster(uoa1, uoa2), data = simdata)
mod1 <- lmitt(y ~ 1, data = simdata, specification = spec, weights = "ate")
mod2 <- lmitt(y ~ as.factor(o), data = simdata, specification = spec, weights = "ate")
### observational study with treatment z assigned row-wise within blocks:
mod3 <- lmitt(y ~ 1, data=simdata, specification=z ~ block(bid), weights="att")
### regression discontinuity study with units of assignment
### given by combinations of uoa1, uoa2:
mod4 <- lmitt(y ~ 1, data = simdata,
```

```
specification = z ~ uoa(uoa1, uoa2) + forcing(force))
```

---

IsoSynth

*Synthetic Regression Discontinuity Data*

---

## Description

The data for this example were randomly simulated using the **synthpop** package in R based on data originally collected by Lindo, Sanders, and Oreopoulos (2010).

## Usage

```
IsoSynth
```

## Format

A data.frame with 40,403 rows and 11 columns.

- R
- lhsgrade\_pct
- nextGPA
- probation\_year1
- totcredits\_year1
- male
- loc\_campus1
- loc\_campus2
- bpl\_north\_america
- english
- age\_at\_entry

## Details

See the "Regression Discontinuity StudySpecifications" vignette on the propertee website for more details on the original data, a link to the code used to generate this synthetic data, and a detailed example.

---

michigan\_school\_pairs *Intervention data from a pair-matched study of schools in Michigan*


---

### Description

Michigan high schools, with a plausible cluster RCT

### Usage

michigan\_school\_pairs

### Format

A data.frame with 14 rows and 13 columns.

- schoolid school id
- blk block
- z treatment variable
- MALE\_G11\_PERC percentage of G11 male students
- FEMALE\_G11\_PERC percentage of G11 female students
- AM\_G11\_PERC percentage of G11 American Indian/Alaska Native students
- ASIAN\_G11\_PERC percentage of G11 Asian students
- HISP\_G11\_PERC percentage of G11 Hispanic students
- BLACK\_G11\_PERC percentage of G11 Black students
- WHITE\_G11\_PERC percentage of G11 White students
- PACIFIC\_G11\_PERC percentage of G11 Hawaiian Native/Pacific Islander students
- TR\_G11\_PERC percentage of G11 Two or More Races students
- G11 Number of G11 students

### Details

Grade 11 demographics for all Michigan high schools in 2013, with mock block and treatment assignments for 14 high schools within a large county in the metro Detroit area. These schools were selected for this demonstration based on their similarity to the 14 high schools from an adjacent Michigan county that participated in the Pane et al (2013) study. As a result, they serve as an example of what one might expect to find as the state-specific school-level subsample in a multi-state paired cluster randomized trial featuring random assignment at the school level.

The mock experimental schools were selected by optimal matching of experimental schools to adjacent county schools, with substitute schools grouped into the same pairs or triples ('fine strata') as were their experimental counterparts. The original pairs and triples had been selected to reduce variation in baseline variables predictive of outcomes, and the blocking structure the substitute sample inherits may be expected to do this as well. The treatment/control distinction is also inherited from the experimental sample, but there is of course no treatment effect within the mock experiment.



The selection of mock experimental schools was based on both demographic and student achievement variables, but the present data frame includes only the demographic variables (as sourced from the Common Core of Data [CCD; U.S. Department of Education]). School average outcomes in student test scores are available separately, from Michigan's Center for Education Performance Information. See the vignette 'Real-data demonstration with a finely stratified cluster RCT and a broader administrative database', available on the package website.

## References

Pane, John F., et al. "Effectiveness of cognitive tutor algebra I at scale." *Educational Evaluation and Policy Analysis* 36.2 (2014): 127-144.

U.S. Department of Education. Public Elementary/Secondary School Universe Survey Data, v.2a. Institute of Education Sciences, National Center for Education Statistics.

## Examples

```
data(michigan_school_pairs)
mi_spec <- rct_spec(z ~ uoa(schoolid)+block(blk),
  data=michigan_school_pairs)
mi_spec
table(is.na(michigan_school_pairs$blk))
specification_table(mi_spec, "block", "treatment")
```

---

rct_spec	<i>Generates a StudySpecification object with the given specifications.</i>
----------	---

---

## Description

Generate a randomized control treatment StudySpecification ([rct\\_spec\(\)](#)), or an observational StudySpecification ([obs\\_spec\(\)](#)), or a regression discontinuity StudySpecification ([rd\\_spec\(\)](#)).

## Usage

```
rct_spec(formula, data, subset = NULL, na.fail = TRUE)

rd_spec(formula, data, subset = NULL, na.fail = TRUE)

obs_spec(formula, data, subset = NULL, na.fail = TRUE)

rct_specification(formula, data, subset = NULL, na.fail = TRUE)

rd_specification(formula, data, subset = NULL, na.fail = TRUE)

obs_specification(formula, data, subset = NULL, na.fail = TRUE)

obsstudy_spec(formula, data, subset = NULL, na.fail = TRUE)

obsstudy_specification(formula, data, subset = NULL, na.fail = TRUE)
```

## Arguments

<code>formula</code>	a formula defining the <code>StudySpecification</code> components. See Details for specification.
<code>data</code>	the data set from which to build the <code>StudySpecification</code> . Note that this data need not be the same as used to estimate the treatment effect; rather the data passed should contain information about the units of treatment assignment (as opposed to the units of analysis).
<code>subset</code>	optional, subset the data before creating the <code>StudySpecification</code> object
<code>na.fail</code>	If TRUE (default), any missing data found in the variables specified in <code>formula</code> (excluding treatment) will trigger an error. If FALSE, non-complete cases will be dropped before the creation of the <code>StudySpecification</code>

## Details

The formula should include exactly one `unit_of_assignment()` to identify the units of assignment (one or more variables). (`uoa`, `cluster`, or `unitid` are synonyms for `unit_of_assignment`; the choice of which has no impact on the analysis. See below for a limited exception in which the `unit_of_assignment` specification may be omitted.) If defining an `rd_spec`, the formula must also include a `forcing()` entry. The formula may optionally include a `block()` as well. Each of these can take in multiple variables, e.g. to pass both a household ID and individual ID as unit of assignment, use `uoa(hhid, iid)` and not `uoa(hhid) + uoa(iid)`.

The treatment variable passed into the left-hand side of formula can either be logical, numeric, or character. If it is anything else, it attempts conversion to one of those types (for example, factor and ordered are converted to numeric if the levels are numeric, otherwise to character). If the treatment is not logical or numeric with only values 0 and 1, in order to generate weights with `ate()` or `ett()`, the dichotomy argument must be used in those functions to identify the treatment and control groups. See `ett()` for more details on specifying a dichotomy.

There are a few aliases for each version.

If the formula excludes a `unit_of_assignment()`, data merges are performed on row order. Such formulas can also be passed as the specification argument to `lmitt()`, and that is their primary intended use case. It is recommended that each formula argument passed to `*_specification()` include a `unit_of_assignment()`, `uoa()` or `cluster()` term identifying the key variable(s) with which `StudySpecification` data is to be merged with analysis data. Exceptions to this rule will be met with a warning. To disable the warning, run `options("propertee_warn_on_no_unit_of_assignment" = FALSE)`.

The units of assignment, blocks, and forcing variables must be numeric or character. If they are otherwise, an attempt is made to cast them into character.

## Value

a `StudySpecification` object of the requested type for use in further analysis.

## Examples

```
data(simdata)
spec <- rct_spec(z ~ unit_of_assignment(uoa1, uoa2) + block(bid),
  data = simdata)
```

```
data(schooldata)
spec <- obs_spec(treatment ~ unit_of_assignment(schoolid) + block(state),
  data = schooldata)
```

---

schooldata

*Student data*


---

## Description

An example of data sets stored at two levels.

## Usage

schooldata

studentdata

## Format

Two `data.frames`, one with school-level data (`schooldata`) including treatment assignment and a second with student-level data (`studentdata`). `schoolata`:

- `schoolid` Unique school ID variable.
- `treatment` Was this school in the intervention group?
- `state` State which the school is in.
- `pct_disadvantage` Percent of student body flagged as "disadvantaged".

`studentdata`:

- `id` Unique student ID.
- `schoolid` Unique school ID variable.
- `grade` Student's grade, 3-5.
- `gpa` Student GPA in prior year.
- `math` Standardized math score (out of 100).

An object of class `data.frame` with 8713 rows and 5 columns.

## Details

In this hypothetical data, schools were randomly assignment to treatment status, but the unit of analysis is students. Thus the two data sets, one encoding school information (including treatment status) and one encoding student information (which does not include treatment status).

**Examples**

```
soec <- obs_spec(treatment ~ uoa(schoolid), data = schooldata)

# Treatment effect
mod1 <- lmitt(math ~ 1, specification = soec, data = studentdata)

# Treatment effect by grade
mod2 <- lmitt(math ~ as.factor(grade), specification = soec, data = studentdata)
```

---

```
show, PreSandwichLayer-method
```

*Show a PreSandwichLayer or SandwichLayer*

---

**Description**

Display information about a PreSandwichLayer or SandwichLayer object

**Usage**

```
## S4 method for signature 'PreSandwichLayer'
show(object)
```

**Arguments**

object                      PreSandwichLayer or SandwichLayer object

**Value**

an invisible copy of object

---

```
show, StudySpecification-method
```

*Show a StudySpecification*

---

**Description**

Display information about a StudySpecification object

**Usage**

```
## S4 method for signature 'StudySpecification'
show(object)
```

**Arguments**

object                      StudySpecification object, usually a result of a call to `rct_spec()`, `obs_spec()`, or `rd_spec()`.

**Value**

object, invisibly.

---

show, teeMod-method	<i>Show a teeMod</i>
---------------------	----------------------

---

**Description**

Display information about a teeMod object

**Usage**

```
## S4 method for signature 'teeMod'
show(object)
```

**Arguments**

object                      teeMod object, usually a result of a call to `lmitt()`.

**Value**

object, invisibly.

---

show,WeightedStudySpecification-method	<i>Show a WeightedStudySpecification</i>
--	--

---

**Description**

Prints out the weights from a WeightedStudySpecification

**Usage**

```
## S4 method for signature 'WeightedStudySpecification'
show(object)
```

**Arguments**

object                      a WeightedStudySpecification object

**Value**

an invisible copy of object

---

simdata

*Simulated data*


---

### Description

Simulated data to use with the **propertee** package with unit of assignment level treatment assignment

### Usage

```
simdata
```

### Format

A data.frame with 100 rows and 7 columns.

- uoa1 First level unit of assignment ID
- uoa2 Second level unit of assignment ID
- bid Block ID
- force Forcing variable
- z Binary treatment indicator
- o 4-level ordered treatment variable
- dose Dose treatment variable
- x Some predictor
- y Some outcome

---

specification\_data\_concordance

*Check for variable agreement within units of assignment*


---

### Description

Useful for debugging purposes to ensure that there is concordance between variables in the StudySpecification and data.

### Usage

```
specification_data_concordance(
  specification,
  data,
  by = NULL,
  warn_on_nonexistence = TRUE
)
```

**Arguments**

specification a StudySpecification object

data a new data set, presumably not the same used to create specification.

by optional; named vector or list connecting names of variables in specification to variables in data. Names represent variables in specification; values represent variables in data. Only needed if variable names differ.

warn\_on\_nonexistence default TRUE. If a variable does not exist in data, should this be flagged? If FALSE, silently move on if a variable doesn't exist in data.

**Details**

Consider the following scenario: A StudySpecification is generated from some dataset, "data1", which includes a block variable "b1". Within each unique unit of assignment/unitid/cluster of "data1", it must be the case that "b1" is constant. (Otherwise the creation of the StudySpecification will fail.)

Next, a model is fit which includes weights generated from the StudySpecification, but on dataset "data2". In "data2", the block variable "b1" also exists, but due to some issue with data cleaning, does not agree with "b1" in "data1".

This could cause errors, either directly (via actual error messages) or simply produce nonsense results. `specification_data_concordance()` is specificationed to help debug these scenarios by providing information on whether variables in both the data used in the creation of specification ("data1" in the above example) and some new dataset, data, ("data2" in the above example) have any inconsistencies.

**Value**

invisibly TRUE if no warnings are produced, FALSE if any warnings are produced.

---

specification_table	<i>Table of elements from a StudySpecification</i>
---------------------	--

---

**Description**

Produces a table (1-dimensional, or 2-dimensional if y is specified) of the elements of the StudySpecification.

**Usage**

```
specification_table(
  specification,
  x,
  y = NULL,
  sort = FALSE,
  decreasing = TRUE,
  use_var_names = FALSE,
```

```

    ...
  )

  stable(
    specification,
    x,
    y = NULL,
    sort = FALSE,
    decreasing = TRUE,
    use_var_names = FALSE,
    ...
  )

```

### Arguments

<code>specification</code>	A <code>StudySpecification</code> object
<code>x</code>	One of "treatment", "unit of assignment", (synonym "uoa"), "block". Abbreviations are accepted. "unit of assignment" can be replaced by "unitid" or "cluster" if the <code>StudySpecification</code> was created with that element.
<code>y</code>	Optionally, another string similar to <code>x</code> . A 1-dimensional table is produced if <code>y</code> is left at its default, <code>NULL</code> .
<code>sort</code>	Ignored if <code>y</code> is not <code>NULL</code> . If <code>FALSE</code> (default), one-way table is sorted according to "names" of levels. If set to <code>TRUE</code> , one-way table is sorted according to values.
<code>decreasing</code>	If <code>sort</code> is <code>TRUE</code> , choose whether to sort descending ( <code>TRUE</code> , default) or ascending ( <code>FALSE</code> ).
<code>use_var_names</code>	If <code>TRUE</code> , name dimensions of table returned by variable names. If <code>FALSE</code> (default), name by their function (e.g. "treatment" or "blocks"). Passing the <code>dnn</code> argument in <code>...</code> (an argument of <code>table()</code> ) overrides whatever is requested here.
<code>...</code>	additional arguments <code>table()</code>

### Value

A table of the requested variables.

### Examples

```

data(simdata)
spec <- obs_spec(z ~ unit_of_assignment(uoa1, uoa2) + block(bid),
  data = simdata)
specification_table(spec, "treatment")
specification_table(spec, "treatment", "block", sort = TRUE, use_var_names = TRUE)

```



---

STARplus

---

*STAR participants plus nonexperimental controls*


---

## Description

Data from Tennessee's Project STAR study. This data frame describes student participants in the Project STAR (Student-Teacher Achievement Ratio) field experiment conducted in Tennessee, USA beginning in the mid-1980s, as well as an external control group consisting of the contemporaneous cohort of students attending a matched sample of Tennessee schools that did not participate in the STAR experiment. Variables are as described in Project STAR data documentation (see references), with five exceptions. Three \*\_at\_entry variables were constructed as follows: grade\_at\_entry indicates the grade of student's first participation, while school\_at\_entry and cond\_at\_entry reflect the school ID and classroom type corresponding to the student's grade at entry to the study. Additionally, read\_yr1 and math\_yr1 capture a student's scaled scores on the Scholastic Assessment Test (SAT) administered to them during their grade\_of\_entry, i.e. their earliest available post-treatment SAT measurements.

## Usage

```
STARplus
```

## Format

A data.frame with 13,382 rows and 56 columns.

- stdnid Student ID
- gender Student gender
- race Student race
- birthmonth Student month of birth
- birthday Student day of birth
- birthyear Student year of birth
- read\_yr1 SAT reading scaled score from grade at which student entered the study
- math\_yr1 SAT math scaled score from grade at which student entered the study
- gktreadss Kindergarten reading scaled score (RCT participants only)
- gktmathss Kindergarten math scaled score (RCT participants only)
- gkltlistss Kindergarten listening scaled score (RCT participants only)
- gkwordskillss Kindergarten word study skills scaled score (RCT participants only)
- g1schid Grade 1 School ID
- g1tchid Grade 1 Teacher ID
- g1classsize Class size of Grade 1
- g1treadss Grade 1 SAT reading scaled score
- g1tmathss Grade 1 SAT math scaled score

- g1tlistss Grade 1 total listening scale score in SAT
- g1wordskillss Grade 1 word study skills scale score in SAT
- g1readbsraw Grade 1 reading raw score in Basic Skills First (BSF) tests
- g1mathbsraw Grade 1 math raw score in BSF
- g1readbsobjpct Grade 1 reading percent objectives mastered in BSF tests
- g1mathbsobjpct Grade 1 math percent objectives mastered in BSF tests
- g2schid Grade 2 School ID
- g2tchid Grade 2 Teacher ID
- g2classsize Class size of Grade 2
- g2treadss Grade 2 total reading scale score in SAT
- g2tmathss Grade 2 total math scale score in SAT
- g2tlistss Grade 2 total listening scale score in SAT
- g2wordskillss Grade 2 word study skills scale score in SAT
- g2readbsraw Grade 2 reading raw score in BSF tests
- g2mathbsraw Grade 2 math raw score in BSF test
- g2readbsobjpct Grade 2 reading percent objectives mastered in BSF tests
- g3schid Grade 3 School ID
- g3tchid Grade 3 Teacher ID
- g3classsize Class size of Grade 3
- g3treadss Grade 3 total reading scale score in SAT
- g3tmathss Grade 3 total math scale score in SAT
- g3langss Grade 3 total language scale score in SAT
- g3tlistss Grade 3 total listening scale score in SAT
- g3socialsciss Grade 3 social science scale score in SAT
- g3spellss Grade 3 spelling scale score in SAT
- g3vocabss Grade 3 vocabulary scale score in SAT
- g3mathcomputss Grade 3 math computation scale score in SAT
- g3mathnumconcss Grade 3 concept of numbers scale score in SAT
- g3mathapplss Grade 3 math applications scale score in SAT
- g3wordskillss Grade 3 word study skills scale score in SAT
- g3readbsraw Grade 3 reading raw score in BSF tests
- g3mathbsraw Grade 3 math raw score in BSF tests
- g3readbsobjpct Grade 3 reading percent objectives mastered in BSF tests
- g3mathbsobjpct Grade 3 math percent objectives mastered in BSF tests
- dob Date of birth (with NAs imputed RCT participant median)
- dobNA Dat of birth not recorded
- grade\_at\_entry Grade at which each student first entered the study
- school\_at\_entry School ID corresponding to the student's grade at entry into the study
- cond\_at\_entry Classroom type corresponding to the student's grade at entry into the study

**Details**

Note: This dataset bears a Creative Commons Zero license (v1.0).

**Source**

[doi:10.7910/DVN/SIWH9F](https://doi.org/10.7910/DVN/SIWH9F)

**References**

C.M. Achilles; Helen Pate Bain; Fred Bellott; Jayne Boyd-Zaharias; Jeremy Finn; John Folger; John Johnston; Elizabeth Word, 2008, "Tennessee's Student Teacher Achievement Ratio (STAR) project", Harvard Dataverse, V1, <https://doi.org/10.7910/DVN/SIWH9F> UNF:3:Ji2Q+9HCCZAbw3csOdMNdA

---

subset,PreSandwichLayer-method  
PreSandwichLayer and SandwichLayer *subsetting*

---

**Description**

Return subset of a PreSandwichLayer or SandwichLayer which meets conditions.

**Usage**

```
## S4 method for signature 'PreSandwichLayer'
subset(x, subset)

## S4 method for signature 'PreSandwichLayer'
x[i]
```

**Arguments**

x	PreSandwichLayer or SandwichLayer object
subset	Logical vector identifying values to keep or drop
i	indices specifying elements to extract or replace. See <code>help("[")</code> for further details.

**Value**

x subset by subset or i

---

```
subset, WeightedStudySpecification-method
      WeightedStudySpecification subsetting
```

---

### Description

Provides functionality to subset the weights of a WeightedStudySpecification object.

### Usage

```
## S4 method for signature 'WeightedStudySpecification'
subset(x, subset)

## S4 method for signature 'WeightedStudySpecification'
x[i]
```

### Arguments

x	WeightedStudySpecification object
subset	Logical vector identifying values to keep or drop
i	indices specifying elements to extract or replace. See <code>help("[")</code> for further details.

### Value

A WeightedStudySpecification object which is a subsetting version of x.

---

```
summary.StudySpecification
      Summarizing StudySpecification objects
```

---

### Description

[summary\(\)](#) method for class StudySpecification.

### Usage

```
## S3 method for class 'StudySpecification'
summary(object, ..., treatment_binary = TRUE)

## S3 method for class 'summary.StudySpecification'
print(x, ..., max_unit_print = 3)
```

Arguments

object	StudySpecification object, usually a result of a call to <a href="#">rct_spec()</a> , <a href="#">obs_spec()</a> , or <a href="#">rd_spec()</a> .
...	Ignored
treatment_binary	Should the treatment be dichotomized if object contains a dichotomy? Ignored if object does not contain a dichotomy.
x	summary.StudySpecification object, usually as a result of a call to <a href="#">summary.StudySpecification()</a>
max_unit_print	Maximum number of treatment levels to print in treatment table

Value

The StudySpecification or summary.StudySpecificationobject, invisibly

---

summary.teeMod	<i>Summarizing teeMod objects</i>
----------------	-----------------------------------

---

Description

[summary\(\)](#) method for class teeMod

Usage

```
## S3 method for class 'teeMod'
summary(object, vcov.type = "HC0", ...)

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

Arguments

object	teeMod object
vcov.type	A string indicating the desired variance estimator. See <a href="#">vcov_tee()</a> for details on accepted types.
...	Additional arguments to <a href="#">vcov_tee()</a> , such as the desired finite sample heteroskedasticity-robust standard error adjustment.
x	summary.teeMod object
digits	the number of significant digits to use when printing.
signif.stars	logical. If 'TRUE', 'significance stars' are printed for each coefficient.

## Details

If a `teeMod` object is fit with a `SandwichLayer` offset, then the usual `stats::summary.lm()` output is enhanced by the use of covariance-adjusted sandwich standard errors, with t-test values recalculated to reflect the new standard errors.

## Value

object of class `summary.teeMod`

---

treatment

*Accessors and Replacers for StudySpecification objects*

---

## Description

Allows access to the elements which define a `StudySpecification`, enabling their extraction or replacement.

## Usage

```
treatment(x, newdata = NULL, dichotomy = NULL, by = NULL, ...)
```

```
## S4 method for signature 'StudySpecification'
```

```
treatment(x, newdata = NULL, dichotomy = NULL, by = NULL, ...)
```

```
treatment(x) <- value
```

```
## S4 replacement method for signature 'StudySpecification'
```

```
treatment(x) <- value
```

```
units_of_assignment(x, newdata = NULL, by = NULL)
```

```
## S4 method for signature 'StudySpecification'
```

```
units_of_assignment(x, newdata = NULL, by = NULL)
```

```
units_of_assignment(x) <- value
```

```
## S4 replacement method for signature 'StudySpecification'
```

```
units_of_assignment(x) <- value
```

```
clusters(x, newdata = NULL, by = NULL)
```

```
## S4 method for signature 'StudySpecification'
```

```
clusters(x, newdata = NULL, by = NULL)
```

```
clusters(x) <- value
```

```
## S4 replacement method for signature 'StudySpecification'
```

```

clusters(x) <- value

unitids(x)

## S4 method for signature 'StudySpecification'
unitids(x)

unitids(x) <- value

## S4 replacement method for signature 'StudySpecification'
unitids(x) <- value

blocks(x, newdata = NULL, by = NULL, ...)

## S4 method for signature 'StudySpecification'
blocks(x, newdata = NULL, by = NULL, ..., implicit = FALSE)

blocks(x) <- value

## S4 replacement method for signature 'StudySpecification'
blocks(x) <- value

has_blocks(x)

forcings(x, newdata = NULL, by = NULL)

## S4 method for signature 'StudySpecification'
forcings(x, newdata = NULL, by = NULL)

forcings(x) <- value

## S4 replacement method for signature 'StudySpecification'
forcings(x) <- value

```

## Arguments

<code>x</code>	a <code>StudySpecification</code> object
<code>newdata</code>	optional; an additional <code>data.frame</code> . If passed, and the unit of assignment variable is found in <code>newdata</code> , then the requested variable type for each unit of <code>newdata</code> is returned. See <code>by</code> argument if the name of the unit of assignment differs.
<code>dichotomy</code>	optional; a formula specifying how to dichotomize a non-binary treatment variable. See the Details section of the <code>ett()</code> or <code>att()</code> help pages for information on specifying this formula
<code>by</code>	optional; named vector or list connecting names of unit of assignment/unitid/cluster variables in <code>x</code> to unit of assignment/unitid/cluster variables in <code>data</code> . Names represent variables in <code>x</code> ; values represent variables in <code>newdata</code> . Only needed if variable names differ.

...	ignored.
value	replacement. Either a vector/matrix of appropriate dimension, or a named data.frame if renaming variable as well. See Details.
implicit	Should a block-less StudySpecification return a constant 1 when extracting blocks?

## Details

For `treatment()`, when argument `binary` is `FALSE`, the treatment variable passed into the `StudySpecification` is returned as a one-column data.frame regardless of whether it is binary or `x` has a dichotomy

If a dichotomy is passed, a binary one-column data.frame will be returned. If not and `binary` is `TRUE`, unless the `StudySpecification` has a binary treatment, `treatment()` will error. If `binary` is "ifany", it will return the original treatment in this case.

The one-column data.frame returned by `treatment()` is named as entered in the `StudySpecification` creation, but if a dichotomy is passed, the column name is `"__z"` to try and avoid any name conflicts.

For the value when using replacers, the replacement must have the same number of rows as the `StudySpecification` (the same number of units of assignment). The number of columns can differ (e.g. if the `StudySpecification` were defined with two variable uniquely identifying blocks, you can replace that with a single variable uniquely identifying blocks, as long as it respects other restrictions.)

If the replacement value is a data.frame, the name of the columns is used as the new variable names. If the replacement is a matrix or vector, the original names are retained. If reducing the number of variables (e.g., moving from two variables uniquely identifying to a single variable), the appropriate number of variable names are retained. If increasing the number of variables, a data.frame with names must be provided.

## Value

data.frame containing requested variable, or an updated `StudySpecification`. `treatment()` works slightly differently, see Details.

## Examples

```
data(simdata)
spec <- obs_spec(z ~ unit_of_assignment(uoa1, uoa2), data = simdata)
blocks(spec) # empty
blocks(spec) <- data.frame(blks = c(1, 1, 2, 2, 3, 3, 4, 4, 5, 5))
blocks(spec)
blocks(spec) <- c(5, 5, 4, 4, 3, 3, 2, 2, 1, 1)
blocks(spec) # notice that variable is not renamed
```



---

unit_of_assignment	<i>Special terms in StudySpecification creation formula</i>
--------------------	---

---

## Description

These are special functions used only in the definition of StudySpecification objects. They identify the units of assignment, blocks and forcing variables. They should never be used outside of the formula argument to obs\_spec, rct\_spec, or rd\_spec.

## Usage

```
unit_of_assignment(...)
```

```
unitid(...)
```

```
cluster(...)
```

```
uoa(...)
```

```
block(...)
```

```
forcing(...)
```

## Arguments

...                    any number of variables of the same length.

## Details

These functions have no use outside of the formula in creating a StudySpecification.

unit\_of\_assignment, uoa, cluster and unitid are synonyms; you must include one and only one in each StudySpecification. The choice of which to use will have no impact on any analysis, only on some output and the name of the stored element in the StudySpecification. Accessors/replacers (units\_of\_assignment, unitids, clusters) respect the choice made at the point of creation of the StudySpecification, and only the appropriate function will work.

See rct\_spec, obs\_spec, or rd\_spec for examples of their usage.

## Value

the variables with appropriate labels. No use outside of their inclusion in the formula argument to obs\_spec, rct\_spec, or rd\_spec

---

var_table	<i>Extract Variable Names from StudySpecification</i>
-----------	---

---

## Description

Methods to extract the variable names to the elements of the structure of the StudySpecification (e.g. treatment, unit of analysis, etc)

## Usage

```
var_table(specification, compress = TRUE, report_all = FALSE)
```

```
var_names(specification, type, implicitBlocks = FALSE)
```

## Arguments

specification	a StudySpecification object
compress	should multiple variables be compressed into a comma-separated string? Default TRUE. If FALSE, multiple columns can be created instead.
report_all	should we report all possible structures even if they don't exist in the StudySpecification? Default FALSE.
type	one of "t", "u", "b", "f"; for "treatment", "unit_of_assignment", "block", and "forcing" respectively
implicitBlocks	If the StudySpecification is created without blocks, setting this to TRUE will return ".blocks_internal" as the variable name corresponding to the blocks.

## Details

When compress is TRUE, the result will always have two columns. When FALSE, the result will have number of columns equal to the largest number of variables in a particular role, plus one. E.g., a call such as `rct_spec(z ~ unitid(a, b, c, d) ...` will have 4+1=5 columns in the output matrix with `compress = FALSE`.

When `report_all` is TRUE, the matrix is guaranteed to have 3 rows (when the specification is an RCT or Obs) or 4 rows (when the specification is a RD), with empty variable entries as appropriate. When FALSE, the matrix will have minimum 2 rows (treatment and unit of assignment/unitid/cluster), with additional rows for blocks and forcing if included in the StudySpecification.

## Value

`var_table` returns the requested table. `var_names` returns a vector of variable names.

## Examples

```
spec <- rct_spec(z ~ uoa(uoa1, uoa2) + block(bid), data = simdata)
var_table(spec)
var_table(spec, compress = FALSE)
var_names(spec, "t")
var_names(spec, "u")
var_names(spec, "b")
```

---

vcov.teeMod

---

*Compute variance-covariance matrix for fitted teeMod model*


---

## Description

An S3method for `stats::vcov` that computes standard errors for `teeMod` models using `vcov_tee()`.

## Usage

```
## S3 method for class 'teeMod'
vcov(object, ...)
```

## Arguments

<code>object</code>	a fitted <code>teeMod</code> model
<code>...</code>	additional arguments to <code>vcov_tee()</code> .

## Details

`vcov.teeMod()` wraps around `vcov_tee()`, so additional arguments passed to `...` will be passed to the `vcov_tee()` call. See documentation for `vcov_tee()` for information about necessary arguments.

## Value

A variance-covariance matrix with row and column entries for the estimated coefficients in `x`, the marginal mean outcome in the control condition, the marginal mean offset in the control condition (if an offset is provided), and if a moderator variable is specified in the formula for `x`, the mean interaction in the control condition of the outcome and offset with the moderator variable

---

`weights, WeightedStudySpecification-method`*Extract Weights from WeightedStudySpecification*

---

**Description**

A `WeightedStudySpecification` object contains a numeric vector with a few additional slots, this extracts only the numeric vector.

**Usage**

```
## S4 method for signature 'WeightedStudySpecification'
weights(object, ...)
```

**Arguments**

<code>object</code>	a <code>WeightedStudySpecification</code> object
<code>...</code>	Ignored

**Value**

A numeric vector of the weights

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