

# Package: hydrorecipes (via r-universe)

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**Type** Package

**Title** Hydrogeology steps

**Version** 0.0.6

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**Description** This package is an implementation of some common steps of the `recipes` package using `R6` classes. It also provides some additional steps that may be useful for the geosciences and signal processing. Two goals of this package are to provide a higher performance package (memory and computation time), and as a learning experience.

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**Encoding** UTF-8

**LazyData** true

**Depends** R (>= 3.5.0), Bessel

**Imports** R6 (>= 2.3.0), Rcpp, data.table, earthtide, collapse, gslns, plotly

**LinkingTo** Rcpp, RcppArmadillo, RcppThread, RcppEigen, splines2, BH

**Suggests** rmarkdown, tinytest, RcppRoll, recipes, tibble, knitr, bench, scales, ggplot2, splines2

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---

areal\_rojstaczer\_semiconfined  
*areal\_rojstaczer\_semiconfined*

---

**Description**

areal\_rojstaczer\_semiconfined

**Usage**

```
areal_rojstaczer_semiconfined(  

  frequency,  

  radius_well,  

  transmissivity,  

  storage_confining,  

  storage_aquifer,  

  diffusivity_confining,  

  diffusivity_vadose,  

  thickness_confining,  

  thickness_vadose,  

  loading_efficiency,  

  attenuation  

)
```

**Arguments**

frequency	the frequency of the response
radius_well	well radius
transmissivity	aquifer transmissivity ( $L^2/t$ )
storage_confining	confining layer storativity ( $L/L$ )
storage_aquifer	aquifer storativity ( $L/L$ )
diffusivity_confining	confining layer diffusivity
diffusivity_vadose	air diffusivity of vadose zone
thickness_confining	confining layer thickness
thickness_vadose	vadose thickness
loading_efficiency	the loading efficiency of the aquifer
attenuation	an attenuation factor

**Value**

complex response vector in frequency domain

---

areal\_rojstaczer\_unconfined  
*areal\_rojstaczer\_unconfined*

---

**Description**

areal\_rojstaczer\_unconfined

**Usage**

```
areal_rojstaczer_unconfined(  
  frequency,  
  radius_well,  
  storage_aquifer,  
  specific_yield,  
  k_vertical,  
  diffusivity_vertical,  
  diffusivity_vadose,  
  thickness_saturated_well,  
  thickness_vadose,
```

```

    thickness_aquifer,
    loading_efficiency,
    attenuation
)

```

### Arguments

`frequency` the frequency of the response

`radius_well` well radius

`storage_aquifer` aquifer storativity (L/L)

`specific_yield` specific yield of unconfined aquifer

`k_vertical` vertical hydraulic conductivity of vadose zone

`diffusivity_vertical` unconfined layer diffusivity

`diffusivity_vadose` air diffusivity of vadose zone

`thickness_saturated_well` saturated well thickness

`thickness_vadose` vadose thickness

`thickness_aquifer` aquifer thickness

`loading_efficiency` the loading efficiency of the aquifer

`attenuation` an attenuation factor

### Value

complex response vector in frequency domain

---

bake

*bake*

---

### Description

Evaluate the steps and store the recipe results

### Usage

```
bake(.rec, data = NULL)
```

**Arguments**

`.rec` the R6 recipe object.

`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 `lm` is called.

**Value**

an updated recipe

**Examples**

```
rec <- recipe(y~x, data = list(x = rnorm(10), y = rnorm(10))) |>
  step_scale(x) |>
  prep() |>
  bake()
```

---

be\_clark\_cpp

*be\_clark\_cpp*

---

**Description**

Clark 1967 solution for calculating barometric efficiency (Algorithm from Batu 1998, pg 76)

**Usage**

```
be_clark_cpp(dep, ind, lag_space, inverse)
```

**Arguments**

`dep` numeric vector of the dependent variable (ie:water level)

`ind` numeric vector of the independent variable (ie:barometric pressure)

`lag_space` integer spacing for lags, useful for higher frequency monitoring

`inverse` logical whether the barometric relationship is inverse

**Value**

barometric efficiency using Clark's method

**Examples**

```
n <- 1000
baro <- sin(seq(0, 2 * pi, length.out = 1000))
wl <- -0.4 * baro + rnorm(1000, sd = 0.02)
be_clark_cpp(wl, baro, lag_space = 1, inverse = TRUE)
```

---

 be\_correct

*be\_correct*


---

### Description

Adjust values based on the barometric efficiency

### Usage

```
be_correct(
  dat,
  dep = "w1",
  ind = "baro",
  be = 0,
  inverse = TRUE,
  known_mean = NULL
)
```

### Arguments

**be** numeric value of the barometric efficiency

**known\_mean** numeric explicitly enter the mean if known. Otherwise estimate from data.

### Value

numeric vector of corrected values

### Examples

```
library(data.table)
baro = rnorm(1000, sd = 0.01) + 9
w1 <- baro * 0.4 + 18
dat <- data.table(baro, w1)
dat$w1 + be_correct(dat, be=0.4, inverse = FALSE)
dat$w1 + be_correct(dat, be=0.4, inverse = FALSE, known_mean = 9)
# should return ~21.6 = 18 + 0.4 * 9
```

---

 be\_visual

*be\_visual*


---

### Description

Generate dataset for comparing barometric efficiency



---

<code>be_visual_data</code>	<i>be_visual_data</i>
-----------------------------	-----------------------

---

## Description

Generate dataset for comparing barometric efficiency

## Usage

```
be_visual_data(
  dat,
  dep = "wl",
  ind = "baro",
  be_tests = seq(0, 1, 0.1),
  inverse = TRUE
)
```

## Arguments

<code>dat</code>	data that has the independent and dependent variables ( <code>data.table</code> )
<code>dep</code>	name of the dependent variable column (character). This is typically the name for the column holding your water level data.
<code>ind</code>	name of the independent variable column (character). This is typically the name for the column holding your barometric pressure data.
<code>be_tests</code>	vector of barometric efficiencies to test (between 0 and 1) (numeric)
<code>inverse</code>	whether the barometric relationship is inverse ( <code>TRUE</code> means that when the barometric pressure goes up the measured water level goes down (vented transducer, depth to water), <code>FALSE</code> means that when the barometric pressure goes up so does the measured pressure (non-vented transducer)) (logical).

## Value

`data.table` of barometric efficiency compensated datasets

## Examples

```
library(data.table)
be <- 0.43
x <- seq(0, 28*pi, pi / (12*12))

baro <- sin(x) + rnorm(length(x), sd = 0.04)
wl <- -sin(x) * be + rnorm(length(x), sd = 0.04)
dat <- data.table(datetime = as.POSIXct(x * 86400 / (2 * pi),
                                       origin = '1970-01-01', tz = 'UTC'),
                 wl = wl, baro = baro)
be_visual_data(dat)
```

---

be_visual_plot	<i>be_visual_plot</i>
----------------	-----------------------

---

### Description

Plot to compare barometric efficiency. Large datasets may take a long time to plot. Sub-sample should be set to TRUE

### Usage

```
be_visual_plot(dat, time = "datetime", subsample = TRUE)
```

### Arguments

dat	data that has the independent and dependent variables (data.table)
time	name of the column containing the time (character)
subsample	should the data be subsampled for plotting? (logical)

### Value

plotly graph for barometric efficiency estimation with Smith method

### Examples

```
library(plotly)
library(data.table)
be <- 0.43
x <- seq(0, 28 * pi, pi / (12 * 12))

baro <- sin(x) + rnorm(length(x), sd = 0.04)
wl <- -sin(x) * be + rnorm(length(x), sd = 0.04)
dat <- data.table(datetime = as.POSIXct(x * 86400 / (2 * pi),
                                     origin = '1970-01-01', tz = 'UTC'),
                 wl = wl, baro = baro)
dat_be <- be_visual_data(dat)
#be_visual_plot(dat_be) #not run
```

---

bessel_k_cplx	<i>bessel_k_cplx</i>
---------------	----------------------

---

### Description

Modified Bessel function of first kind order 1

**Usage**

```
bessel_k_cplx(x, nu, expon_scaled, n_seq)
```

**Arguments**

<code>x</code>	numeric value to evaluate
<code>nu</code>	numeric value to evaluate
<code>expon_scaled</code>	boolean value to evaluate
<code>n_seq</code>	nseq value to evaluate

**Value**

bessel function result

---

bouwer

*bouwer*

---

**Description**

Keeth elementary dataset 1997-08-14 (DATA sheet Slug\_Bouwer-Rice.xls) <https://pubs.usgs.gov/of/2002/ofr02197Rice.xls>

**Usage**

```
data(bouwer)
```

**Format**

data.table

**Examples**

```
data(bouwer)
```

---

bouwer_abc	<i>bouwer_abc</i>
------------	-------------------

---

**Description**

a, b and c values for Bouwer and Rice, 1976 solution, for testing <https://pubs.usgs.gov/of/2002/ofr02197/spreadsheet/Rice.xls>

**Usage**

```
data(bouwer_abc)
```

**Format**

```
data.table
```

**Examples**

```
data(bouwer_abc)
```

---

bouwer_rice	<i>Calculate transmissivity with Bouwer-Rice solution</i>
-------------	---

---

**Description**

Calculate transmissivity with Bouwer-Rice solution

**Usage**

```
bouwer_rice(time, drawdown, radius_screen, radius_casing, Le, Lw, H)
```

**Arguments**

<code>time</code>	the elapsed time
<code>drawdown</code>	the drawdown
<code>radius_screen</code>	radius of the screen
<code>radius_casing</code>	radius of the casing where the water level is
<code>Le</code>	Effective screen length
<code>Lw</code>	height of water from bottom of well
<code>H</code>	height from bottom of aquifer

**Value**

transmissivity from bouwer\_rice

---

`bouwer_rice_abc`      *Calculate equations 4 and 5 from bouwer, 1989*

---

**Description**

Calculate equations 4 and 5 from bouwer, 1989

**Usage**

```
bouwer_rice_abc(rw, Le, Lw, H)
```

**Arguments**

<code>rw</code>	radius of well
<code>Le</code>	Effective screen length
<code>Lw</code>	height of water from bottom of well
<code>H</code>	height from bottom of aquifer

**Value**

$\ln(R_e/rw)$

---

`convert_for_rojstaczer`  
*convert\_for\_rojstaczer*

---

**Description**

`convert_for_rojstaczer`

**Usage**

```
convert_for_rojstaczer(tf)
```

**Arguments**

<code>tf</code>	the transfer function
-----------------	-----------------------

**Value**

conjugate of tf

---

convert_le_to_be	<i>convert_le_to_be</i>
------------------	-------------------------

---

**Description**

convert\_le\_to\_be

**Usage**

```
convert_le_to_be(tf)
```

**Arguments**

**tf**                    complex transfer function

**Value**

the loading efficiency

---

convolve_filter	<i>convolve_filter</i>
-----------------	------------------------

---

**Description**

convolution of vector with matrix

**Usage**

```
convolve_filter(x, y, remove_partial, reverse)
```

**Arguments**

**x**                    vector to convolve with y (numeric vector)

**y**                    numeric matrix to convolve with x (column by column convolution) (numeric matrix)

**remove\_partial**      keep the end values or fill with NA (boolean)

**reverse**             should x be reversed before convolution (boolean)

**Value**

numeric matrix of convolved values

## Examples

```
a <- convolve_filter(x = 1:100,
                    y = c(1:10, rep(0, 90)),
                    remove_partial = FALSE,
                    reverse = TRUE)

b <- stats::convolve(1:100, rev(1:10), type = 'filter')
```

---

convolve_matrix	<i>convolve_matrix</i>
-----------------	------------------------

---

## Description

convolution of vector with matrix

## Usage

```
convolve_matrix(x, y, remove_partial, reverse)
```

## Arguments

<code>x</code>	vector to convolve with y (numeric vector)
<code>y</code>	numeric matrix to convolve with x (column by column convolution) (numeric matrix)
<code>remove_partial</code>	keep the end values or fill with NA (boolean)
<code>reverse</code>	should x be reversed before convolution (boolean)

## Value

numeric matrix of convolved values

## Examples

```
a <- convolve_matrix(x = 1:100,
                    y = as.matrix(1:10),
                    remove_partial = FALSE,
                    reverse = TRUE)

b <- stats::convolve(1:100, rev(1:10), type = 'filter')
```

---

`distributed_lag_list` *distributed\_lag\_list*

---

## Description

Create distributed lag terms

## Usage

```
distributed_lag_list(
  x,
  n_lag,
  lag_max,
  df,
  degree,
  internal_knots,
  boundary_knots,
  complete_basis,
  periodic,
  derivs,
  integral
)
```

## Arguments

<code>x</code>	numeric vector to lag
<code>n_lag</code>	number of lag terms
<code>lag_max</code>	integer the maximum lag
<code>df</code>	Degree of freedom that equals to the column number of the returned matrix. One can specify <code>df</code> rather than <code>knots</code> , then the function chooses <code>df - degree - as.integer(intercept)</code> internal knots at suitable quantiles of <code>x</code> ignoring missing values and those <code>x</code> outside of the boundary. For periodic splines, <code>df - as.integer(intercept)</code> internal knots will be chosen at suitable quantiles of <code>x</code> relative to the beginning of the cyclic intervals they belong to (see Examples) and the number of internal knots must be greater or equal to the specified <code>degree - 1</code> . If internal knots are specified via <code>knots</code> , the specified <code>df</code> will be ignored.
<code>degree</code>	A nonnegative integer specifying the degree of the piecewise polynomial. The default value is 3 for cubic splines. Zero degree is allowed for piecewise constant basis functions.
<code>internal_knots</code>	location of internal knots
<code>boundary_knots</code>	location of boundary knots
<code>complete_basis</code>	logical intercept?

<b>periodic</b>	A logical value. If <b>TRUE</b> , the periodic splines will be returned. The default value is <b>FALSE</b> .
<b>derivs</b>	A nonnegative integer specifying the order of derivatives of splines basis function. The default value is 0.
<b>integral</b>	A logical value. If <b>TRUE</b> , the corresponding integrals of spline basis functions will be returned. The default value is <b>FALSE</b> . For periodic splines, the integral of each basis is integrated from the left boundary knot.

**Value**

List of distributed lags

---

<code>get_formula_vars</code>	<i>get_formula_vars</i>
-------------------------------	-------------------------

---

**Description**

`get_formula_vars`

**Usage**

```
get_formula_vars(formula, data)
```

**Arguments**

<b>formula</b>	an object of class " <b>formula</b> " (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'.
<b>data</b>	an optional data frame, list or environment (or object coercible by <code>as.data.frame</code> to a data frame) containing the variables in the model. If not found in <b>data</b> , the variables are taken from <code>environment(formula)</code> , typically the environment from which <code>lm</code> is called.

**Value**

character vector of variable names

---

grf_grid	<i>grf_grid</i>
----------	-----------------

---

## Description

Parallel convolution of GRF well function and flow rates in the time domain. Time series needs to be regularly spaced and so are the flow rates. Some performance gains can be achieved if the number of flow rate does not change for each time.

## Usage

```
grf_grid(  
  grid,  
  well_locations,  
  flow_rate,  
  time,  
  specific_storage,  
  hydraulic_conductivity,  
  thickness,  
  flow_dimension  
)
```

## Arguments

<code>grid</code>	locations of grid points (x,y)
<code>well_locations</code>	locations of wells (x,y)
<code>flow_rate</code>	well flow rates
<code>time</code>	prediction times
<code>specific_storage</code>	aquifer specific storage
<code>hydraulic_conductivity</code>	aquifer hydraulic conductivity
<code>thickness</code>	aquifer thickness
<code>flow_dimension</code>	flow dimension

## Value

this solution for multiple pumping scenario

---

<code>grf_time</code>	<i>grf_time</i>
-----------------------	-----------------

---

### Description

Convolution of GRF well function and flow rates in the time domain. Time series needs to be regularly spaced and so are the flow rates. Some performance gains can be achieved if the number of flow rate does not change for each time.

### Usage

```
grf_time(
  radius,
  specific_storage,
  hydraulic_conductivity,
  thickness,
  time,
  flow_rate,
  flow_dimension
)
```

### Arguments

<code>radius</code>	distance to monitoring interval
<code>specific_storage</code>	aquifer storativity
<code>hydraulic_conductivity</code>	aquifer hydraulic conductivity
<code>thickness</code>	aquifer thickness
<code>time</code>	prediction times
<code>flow_rate</code>	well flow rates
<code>flow_dimension</code>	flow dimension

### Value

this solution for multiple pumping scenario

---

hantush_jacob	<i>hantush_jacob</i>
---------------	----------------------

---

### Description

Convolution of hantush well function and flow rates in the time domain. Time series needs to be regularly spaced.

### Usage

```
hantush_jacob(
  time,
  flow_rate,
  radius,
  storativity,
  transmissivity,
  leakage,
  precision
)
```

### Arguments

time	prediction times
flow_rate	well flow rates
radius	distance to monitoring interval
storativity	aquifer storativity
transmissivity	aquifer transmissivity
leakage	hantush leakage
precision	how precise should the solution be. More is more precise but slower.

### Value

hantush jacob solution for multiple pumping scenario

---

hantush_well	<i>hantush_well</i>
--------------	---------------------

---

### Description

Result of the hantush well function

Prodanoff, J.H.A., Mansur, W.J. and Mascarenhas, F.C.B., 2006. Numerical evaluation of Theis and Hantush-Jacob well functions. *Journal of hydrology*, 318(1-4), pp.173-183.

**Usage**

```
hantush_well(u, b, precision)
```

**Arguments**

**u** value of the Theis u  
**b** the leakance  
**precision** how precise should the solution be. More is more precise but slower.

**Value**

hantush well function

---

<code>harmonic_list</code>	<i>harmonic_list</i>
----------------------------	----------------------

---

**Description**

Create sin and cosine terms for harmonic analysis

**Usage**

```
harmonic_list(time, frequency, start, cycle_size)
```

**Arguments**

**time** numeric vector of times  
**frequency** numeric vector of frequencies  
**start** time the cycle starts  
**cycle\_size** size of the cycle in number of measurements

**Value**

List of cosines and sines

---

hsieh\_1987\_fig\_2\_3     *Hsieh (1987) Figures 2 and 3 Digitized*

---

**Description**

Amplitude and phase response as a function of S.

**Usage**

```
hsieh_1987_fig_2_3
```

**Format**

A `data.table` The columns are:

`dimensionless_frequency` dimensionless frequency

`response_gain` and phase of response

`S` storativity

`variable` gain or phase

**Examples**

```
utils::data(hsieh_1987_fig_2_3)
```

---

hussein\_gain     *hussein\_gain*

---

**Description**

Hussein 2013, Figure 5 gain

**Usage**

```
data(hussein_gain)
```

**Format**

`data.table`

**Examples**

```
data(hussein_gain)
```

---

hussein_phase	<i>hussein_phase</i>
---------------	----------------------

---

**Description**

Hussein 2013, Figure 5 phase

**Usage**

```
data(hussein_phase)
```

**Format**

data.table

**Examples**

```
data(hussein_phase)
```

---

kdr	<i>kdr</i>
-----	------------

---

**Description**

measured drawdown from kruseman and de ridder 2000 (figure 3.3). Oude Korendijk pumping test

**Usage**

```
data(kdr)
```

**Format**

data.table

**Examples**

```
data(kdr)
```

---

<code>kelvin</code>	<i>kelvin Kelvin functions of the second kind <math>ker</math> and <math>kei</math> and order 0 to 1.</i>
---------------------	---

---

**Description**

kelvin Kelvin functions of the second kind  $ker$  and  $kei$  and order 0 to 1.

**Usage**

```
kelvin(z, nSeq = 2)
```

**Arguments**

<code>z</code>	value to evaluate the kelvin functions
<code>nSeq</code>	positive integer; if $> 1$ , computes the result for a whole <i>sequence</i> of <code>nu</code> values; if <code>nu</code> $\geq 0$ , <code>nu</code> , <code>nu+1</code> , ..., <code>nu+nSeq-1</code> , if <code>nu</code> $< 0$ , <code>nu</code> , <code>nu-1</code> , ..., <code>nu-nSeq+1</code> .

**Value**

data.table of real and imaginary kelvin functions

---

<code>kennel_2020</code>	<i>kennel_2020</i>
--------------------------	--------------------

---

**Description**

SSFL 2016-08-18 00:00:00 to 2016-10-13 12:00:00 barometric pressure and water pressure data from the same hole. 1 minute interval, pressures in dbar, RBRsoloD 20dbar

**Usage**

```
data(kennel_2020)
```

**Format**

data.table

**Examples**

```
data(kennel_2020)
```

---

lag_list	<i>lag_list</i>
----------	-----------------

---

**Description**

Create lagged terms

**Usage**

```
lag_list(x, lags, n_subset, n_shift)
```

**Arguments**

<code>x</code>	numeric vector - variable to lag
<code>lags</code>	integer vector - amount to lag
<code>n_subset</code>	take every <code>n_subset</code> rows
<code>n_shift</code>	shift values from starting on first row. Should be less than <code>n_subset</code>

**Value**

List of lagged terms

---

liu_1989_fig_8	<i>Liu (1989) Figure 8 Digitized</i>
----------------	--------------------------------------

---

**Description**

Amplitude comparison of Cooper and Liu.

**Usage**

```
liu_1989_fig_8
```

**Format**

A `data.table` The columns are:

<code>period</code>	dimensionless frequency
<code>response</code>	gain of response
<code>aquifer_thickness</code>	thickness of the aquifer (m)
<code>method</code>	cooper or liu method

**Examples**

```
utils::data(liu_1989_fig_8)
```

---

log_lags	<i>log_lags</i>
----------	-----------------

---

**Description**

Generate logarithmically spaced lags

**Usage**

```
log_lags(n, lag_max)
```

**Arguments**

n	integer number of lag terms
lag_max	integer the maximum lag

**Value**

vector of logarithmically spaced lags

---

pad_num	<i>pad_num</i>
---------	----------------

---

**Description**

This function creates a sequence of numbers with 0 padding based on the series length.

**Usage**

```
pad_num(n, pad = "0")
```

**Arguments**

n	the number of columns
pad	the character to use for the padding.

**Value**

a character string padded by "0"

---

plate	<i>plate</i>
-------	--------------

---

## Description

Get the results from the recipe. If the recipe hasn't been prepped and baked, this will do those steps and return the result.

## Usage

```
plate(.rec, type = "dt", ...)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>type</code>	the return type for the recipe (dt = 'data.table', df = 'data.frame', tbl = 'tibble', list = 'list', m = 'matrix')
<code>...</code>	additional arguments

## Value

an updated recipe

## Examples

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_scale(x) |>
  prep() |>
  bake() |>
  plate()

rec <- recipe(y~x, data = dat) |>
  step_scale(x) |>
  plate()
```

---

prep	<i>prep</i>
------	-------------

---

### Description

prep a recipe

### Usage

```
prep(.rec, retain = TRUE)
```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>retain</code>	logical - currently not implemented

### Value

an updated recipe

### Examples

```
rec <- recipe(y~x, data = list(x = rnorm(10), y = rnorm(10))) |>
  step_scale(x) |>
  prep()
```

---

recipe	<i>Create a new R6 recipe. This is analogous to the the list structure that the <i>*recipes*</i> package uses.</i>
--------	--

---

### Description

Create a new recipe object.

### Usage

```
recipe(formula, data, ...)
```

### Arguments

<code>formula</code>	The model formula. It cannot contain operations.
<code>data</code>	list, data.frame, data.table, tibble of data. They will all be treated as lists.
<code>...</code>	additional arguments to pass to Recipe\$new(). This is currently not used.

**Value**

A new R6 'Recipe' object.

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))  
rec <- recipe(y~x, data = dat)
```

---

rojstaczer\_1988a\_fig\_3

*Rojstaczer (1988a) Figure 3 Digitized*

---

**Description**

Barometric efficiency and phase as a function of dimensionless frequency

**Usage**

```
rojstaczer_1988a_fig_3
```

**Format**

A `data.table` The columns are:

`W` dimensionless frequency

`response` gain and phase of response

`variable` gain or phase

**Examples**

```
utils::data(rojstaczer_1988a_fig_3)
```

---

rojstaczer\_1988b\_fig\_3

*Rojstaczer (1988b) Figure 3 Digitized*

---

**Description**

Barometric efficiency and phase as a function of  $Q/W$ . Static confined barometric efficiency is 0.5 and  $R \ll Q$ .  $S$  and  $S'$  are 0.0001.

**Usage**

```
rojstaczer_1988b_fig_3
```

**Format**

A `data.table` The columns are:

`W` dimensionless frequency

`response` gain and phase of response

`Q_div_W` ratio comparing water table drainage and borehole storage

`variable` gain or phase

**Examples**

```
utils::data(rojstaczer_1988b_fig_3)
```

---

`rojstaczer_1988b_fig_5`

*Rojstaczer (1988b) Figure 5 Digitized*

---

**Description**

Amplitude and phase response as a function of  $S$ .

**Usage**

```
rojstaczer_1988b_fig_5
```

**Format**

A `data.table` The columns are:

`W` dimensionless frequency

`response` gain and phase of response

`S` storativity

`variable` gain or phase

**Examples**

```
utils::data(rojstaczer_1988b_fig_5)
```

---

```
rojstaczer_1988b_fig_6
```

*Rojstaczer (1988b) Figure 6 Digitized*

---

### Description

Amplitude and phase response as a function of  $R\_div\_Q$  and dimensionless frequency.

### Usage

```
rojstaczer_1988b_fig_6
```

### Format

A `data.table` The columns are:

`dimensionless_frequency` dimensionless frequency

`response` gain and phase of response

`R_div_Q` R div Q

`variable` gain or phase

### Examples

```
utils::data(rojstaczer_1988b_fig_6)
```

---

```
rojstaczer_1990_fig_2
```

*Rojstaczer and Riley (1990) Figure 2 Digitized*

---

### Description

Amplitude of water table response to Earth tides as a function of  $\Omega'$

### Usage

```
rojstaczer_1990_fig_2
```

### Format

A `data.table` The columns are:

`ohm` water table parameter

`response` amplitude of the response

### Examples

```
utils::data(rojstaczer_1990_fig_2)
```

```
plot(response~ohm, rojstaczer_1990_fig_2, log = 'x')
```

---

`rojstaczer_1990_fig_3`*Rojstaczer and Riley (1990) Figure 3 Digitized*

---

**Description**

Areal strain sensitivity and phase of phreatic well to Earth tides as a function of dimensionless frequency  $Qu$  and partial penetration. Strain sensitivity is 0.5 (I think this is wrong in the paper where it is listed as 0.05)

**Usage**`rojstaczer_1990_fig_3`**Format**

A `data.table` The columns are:

`Qu` dimensionless frequency

`response` gain and phase of response

`b_div_d` partial penetration saturated thickness / aquifer thickness

`variable` gain or phase

**Examples**`utils::data(rojstaczer_1990_fig_3)`

---

`rojstaczer_1990_fig_4`*Rojstaczer and Riley (1990) Figure 4 Digitized*

---

**Description**

Barometric efficiency and phase of phreatic well to atmospheric loading as a function of dimensionless frequency  $Qu$  and  $R/Qu$  and fully penetrating. Static barometric efficiency is 0.5.

**Usage**`rojstaczer_1990_fig_4`

**Format**

A `data.table` The columns are:

`Qu` dimensionless frequency

`response` gain and phase of response

`R_div_Qu` ratio comparing water table drainage and unsaturated zone influences

`variable` gain or phase

**Examples**

```
utils::data(rojstaczer_1990_fig_4)
```

---

<code>step_add_noise</code>	<code>step_add_noise</code>
-----------------------------	-----------------------------

---

**Description**

Add noise.

**Usage**

```
step_add_noise(
  .rec,
  terms,
  sd = 1,
  mean = 0,
  fun = rnorm,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>sd</code>	the standard deviation of the noise to add
<code>mean</code>	the mean of the noise to add
<code>fun</code>	the random noise generating function
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

add noise to a variable

**Examples**

```
dat <- data.frame(x = rnorm(100), y = rnorm(100))
rec <- recipe(y~x, data = dat) |>
  step_add_noise(x) |> plate()
```

---

<code>step_add_vars</code>	<i>step_add_vars</i>
----------------------------	----------------------

---

**Description**

Add a variable from the initial (template) data not included in the recipe creation.

**Usage**

```
step_add_vars(.rec, terms, role = "predictor", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

recipe with added variables

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10), z = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_add_vars(z) |>
  plate()
```

---

```
step_aquifer_constant_drawdown
  step_aquifer_constant_drawdown
```

---

## Description

Estimates the flows of a constant drawdown test using Jacob-Lohman 1952.

## Usage

```
step_aquifer_constant_drawdown(
  .rec,
  time,
  drawdown = 1,
  thickness = 1,
  radius_well = 0.15,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  n_terms = 16L,
  role = "predictor",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>drawdown</code>	drawdown at the well (L)
<code>thickness</code>	the aquifer thickness (L)
<code>radius_well</code>	the radius of the well (L)
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>n_terms</code>	number of terms for laplace solution inversion
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

## References

Jacob, C.E. and S.W. Lohman, 1952. Nonsteady flow to a well of constant drawdown in an extensive aquifer, Trans. Am. Geophys. Union, vol. 33, pp. 559-569.

**See Also**

Other aquifer: [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

**Examples**

```
time <- 10^seq(-5, 2, 0.1)
form <- formula(time~.)
dat <- data.frame(time = time)

jl = recipe(formula = form, data = dat) |>
  step_aquifer_constant_drawdown(time = time,
                                drawdown = 10,
                                thickness = 10,
                                radius_well = 0.15,
                                specific_storage = 1e-6,
                                hydraulic_conductivity = 1,
                                n_terms = 12L) |>
  plate()
```

---

step\_aquifer\_grf      *step\_aquifer\_grf*

---

**Description**

Generates the drawdown using the Generalized Radial Flow (GRF) model. This method defaults to a fast FFT convolution so many rates can be included, but requires a regular time series.

**Usage**

```
step_aquifer_grf(
  .rec,
  time,
  flow_rate,
  thickness = 1,
  radius = 100,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  flow_dimension = 2,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )
<code>thickness</code>	the aquifer thickness (L)
<code>radius</code>	the distance to the observation location (L)
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>flow_dimension</code>	aquifer flow dimension (1 = linear, 2 = radial, 3 = spherical)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

The drawdown using the GRF model

**References**

Barker, J.A., A generalized radial flow model for hydraulic tests in fractured rock. *Water Resour. Res.*, 24 (1988), pp. 1796-1804, 10.1029/WR024i010p01796

**See Also**

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

**Examples**

```
time <- 1:2000
flow_rate <- c(rep(0.001, 500),
               rep(0.002, 500),
               rep(0.0, 1000))

dat <- data.frame(time, flow_rate)

# radial (flow_dimension = 2 Theis)
dd_rad <- recipe(time~flow_rate, dat) |>
  step_aquifer_grf(time = time,
                  flow_rate = flow_rate,
                  thickness = 1.0,
                  radius = 20,
                  specific_storage = 1e-5,
                  hydraulic_conductivity = 1e-3,
                  flow_dimension = 2) |>
  plate()
```

```
# linear (flow_dimension = 1)

dd_lin <- recipe(time~flow_rate, dat) |>
  step_aquifer_grf(time = time,
                   flow_rate = flow_rate,
                   thickness = 1.0,
                   radius = 20,
                   specific_storage = 1e-5,
                   hydraulic_conductivity = 1e-3,
                   flow_dimension = 1) |>
  plate()

# spherical (flow_dimension = 3)
dd_sph <- recipe(time~flow_rate, dat) |>
  step_aquifer_grf(time = time,
                   flow_rate = flow_rate,
                   thickness = 1.0,
                   radius = 20,
                   specific_storage = 1e-5,
                   hydraulic_conductivity = 1e-3,
                   flow_dimension = 3) |>
  plate()
```

---

step\_aquifer\_leaky    *step\_aquifer\_leaky*

---

## Description

hantush\_jacob 1955 solution for a leaky aquifer

## Usage

```
step_aquifer_leaky(
  .rec,
  time,
  flow_rate,
  thickness = 1,
  leakage = 100,
  radius = 100,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  precision = 1e-10,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )
<code>thickness</code>	the aquifer thickness (L)
<code>leakage</code>	the leakage defined by hantush (smaller indicates more leaky)
<code>radius</code>	the distance to the observation location (L)
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>precision</code>	the precision of the solution (default 1e-10)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

The drawdown using the Hantush and Jacob 1955 model

**References**

Prodanoff, J.H.A., Mansur, W.J. and Mascarenhas, F.C.B., 2006. Numerical evaluation of Theis and Hantush-Jacob well functions. *Journal of hydrology*, 318(1-4), pp.173-183. eq: 10, 11, 12

Hantush, M.S. and C.E. Jacob, 1955. Non-steady radial flow in an infinite leaky aquifer, *Am. Geophys. Union Trans.*, vol. 36, no. 1, pp. 95-100.

**See Also**

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

**Examples**

```
time <- 1:2000
flow_rate <- c(rep(0.001, 500),
               rep(0.002, 500),
               rep(0.0, 1000))

# high
dat <- data.frame(time = 1:2000, flow_rate = flow_rate)
hj_100 <- recipe(flow_rate~time, dat) |>
  step_aquifer_leaky(time,
                    flow_rate,
                    leakage = 100,
                    radius = 100,
                    storativity = 1e-6,
```

```

                                transmissivity = 1e-4) |>
  plate()

# medium
hj_200 <- recipe(flow_rate~time, dat) |>
  step_aquifer_leaky(time,
                    flow_rate,
                    leakage = 200,
                    radius = 100,
                    storativity = 1e-6,
                    transmissivity = 1e-4) |>
  plate()

# low
hj_1000 <- recipe(flow_rate~time, dat) |>
  step_aquifer_leaky(time,
                    flow_rate,
                    leakage = 1000,
                    radius = 100,
                    storativity = 1e-6,
                    transmissivity = 1e-4) |>
  plate()

```

---

```
step_aquifer_patch  step_aquifer_patch
```

---

## Description

barker\_herbert 1982 solution for radial patches.

## Usage

```

step_aquifer_patch(
  .rec,
  time,
  flow_rate = 0.01,
  thickness = 1,
  radius = 200,
  radius_patch = 100,
  specific_storage_inner = 1e-06,
  specific_storage_outer = 1e-05,
  hydraulic_conductivity_inner = 1e-04,
  hydraulic_conductivity_outer = 1e-06,
  n_stehfest = 12L,
  role = "predictor",
  ...
)

```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )
<code>thickness</code>	the aquifer thickness (L)
<code>radius</code>	the distance to the observation location (L)
<code>radius_patch</code>	the radius of the cylindrical patch (L)
<code>specific_storage_inner</code>	specific storage of inner patch (L/L)
<code>specific_storage_outer</code>	specific storage of outer patch (L/L)
<code>hydraulic_conductivity_inner</code>	the hydraulic conductivity of the inner patch (L/t)
<code>hydraulic_conductivity_outer</code>	the hydraulic conductivity of the outer patch (L/t)
<code>n_stehfest</code>	integer number of terms to use in Stehfest method (typically < 18)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

The drawdown using the Theis model

**References**

- Barker, J.A., and R. Herbert, 1982: Pumping tests in patchy aquifers, Ground Water, vol. 20, No. 2, pp. 150-155.
- Butler, J.J., 1988: Pumping tests in nonuniform aquifers – The radially symmetric case, Journal of Hydrology, Vol. 101, pp. 15-30.

**See Also**

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

**Examples**

```
dat <- data.frame(time = as.numeric(1:100))
formula <- as.formula(time~.)

frec = recipe(formula = formula, data = dat) |>
  step_aquifer_patch(time = time, flow_rate = 0.01) |>
  plate()
```

---

```
step_aquifer_theis  step_aquifer_theis
```

---

## Description

Radial Flow (GRF) model with flow dimension equal to 2. This method defaults to a fast FFT convolution so many rates can be included, but requires a regular time series.

## Usage

```
step_aquifer_theis(
  .rec,
  time,
  flow_rate,
  thickness = 1,
  radius = 100,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  role = "predictor",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )
<code>thickness</code>	the aquifer thickness (L)
<code>radius</code>	the distance to the observation location (L)
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

The drawdown using the Theis model

## References

Barker, J.A., A generalized radial flow model for hydraulic tests in fractured rock. *Water Resour. Res.*, 24 (1988), pp. 1796-1804, 10.1029/WR024i010p01796

Theis, C.V., 1935: The relation between the lowering of the piezometric surface and the rate and duration of discharge of a well using ground-water storage, *Transactions of the American Geophysical Union*, 16th Annual Meeting, Part 2, pp. 519-524.

**See Also**

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

**Examples**

```
dat <- data.frame(x = as.numeric(1:100),
                 y = rep(0.01, 100))
formula <- as.formula(y~x)

frec = recipe(formula = formula, data = dat) |>
  step_aquifer_theis(time = x, flow_rate = y) |>
  prep() |>
  bake()
```

---

```
step_aquifer_theis_aniso
  step_aquifer_theis_aniso
```

---

**Description**

Generates the drawdown using the Papadopoulos 1965 model. Papadopoulos, I.S., 1965. Nonsteady flow to a well in an infinite anisotropic aquifer.

**Usage**

```
step_aquifer_theis_aniso(
  .rec,
  time,
  flow_rate,
  thickness = 1,
  distance_x = 100,
  distance_y = 100,
  specific_storage = 1e-06,
  hydraulic_conductivity_major = 1e-04,
  hydraulic_conductivity_minor = 1e-05,
  major_axis_angle = 0,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )

<code>thickness</code>	the aquifer thickness (L)
<code>distance_x</code>	distance in the x direction
<code>distance_y</code>	distance in the y direction
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity_major</code>	hydraulic conductivity in the major principal direction
<code>hydraulic_conductivity_minor</code>	hydraulic conductivity in the minor principal direction
<code>major_axis_angle</code>	the orientation of the major principal axis (angle between the x axis and major axis)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

### Value

The drawdown using the Papadopoulos 1965 model

### References

Papadopoulos, I.S., 1965. Nonsteady flow to a well in an infinite anisotropic aquifer. Heilweil, V.M. and Hsieh, P.A., 2006. Determining anisotropic transmissivity using a simplified Papadopoulos method. *Groundwater*, 44(5), pp.749-753.

### See Also

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_wellbore\\_storage\(\)](#)

### Examples

```
dat <- data.frame(x = as.numeric(1:100),
                 y = rep(0.01, 100))
formula <- as.formula(y~x)

frec = recipe(formula = formula, data = dat) |>
  step_aquifer_theis_aniso(time = x, flow_rate = y) |>
  prep() |>
  bake()
```

---

```
step_aquifer_wellbore_storage
  step_aquifer_patch
```

---

## Description

Papadopoulos-Cooper 1967 solution for wellbore storage.

## Usage

```
step_aquifer_wellbore_storage(
  .rec,
  time,
  flow_rate = 1,
  radius = 0.15,
  radius_casing = 0.15,
  radius_well = 0.15,
  thickness = 1,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  n_terms = 12L,
  role = "predictor",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>flow_rate</code>	the flow rate from the well ( $L^3/t$ )
<code>radius</code>	the distance to the observation location (L)
<code>radius_casing</code>	radius of casing in the interval over which the water level declines (L)
<code>radius_well</code>	effective radius of well screen or open hole (L)
<code>thickness</code>	the aquifer thickness (L)
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>n_terms</code>	number of terms for laplace solution inversion
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

The drawdown using the Papadopoulos-Cooper model

## References

Papadopoulos, I.S. and H.H. Cooper, 1967. Drawdown in a well of large diameter, Water Resources Research, vol. 3, no. 1, pp. 241-244.

## See Also

Other aquifer: [step\\_aquifer\\_constant\\_drawdown\(\)](#), [step\\_aquifer\\_grf\(\)](#), [step\\_aquifer\\_leaky\(\)](#), [step\\_aquifer\\_patch\(\)](#), [step\\_aquifer\\_theis\(\)](#), [step\\_aquifer\\_theis\\_aniso\(\)](#)

## Examples

```
dat <- data.frame(x = 10^seq(-5, 2, length.out = 100),
                 y = rep(0.01, 100))
formula <- as.formula(y~x)

frec = recipe(formula = formula, data = dat) |>
  step_aquifer_wellbore_storage(time = x,
                               flow_rate = 10.0,
                               hydraulic_conductivity = 10.0,
                               specific_storage = 1e-4) |>
  prep() |>
  bake()
```

---

step_baro_clark	<i>step_baro_clark</i>
-----------------	------------------------

---

## Description

Clark 1967 solution for calculating barometric efficiency (Algorithm from Batu 1998, pg 76)

## Usage

```
step_baro_clark(
  .rec,
  water_level,
  barometric_pressure,
  lag_space = 1L,
  inverse = FALSE,
  role = "augment",
  ...
)
```

**Arguments**

`.rec` the R6 recipe object.

`water_level` numeric vector of the dependent variable (ie:water level)

`barometric_pressure` numeric vector of the independent variable (ie:barometric pressure)

`lag_space` integer spacing for lags, useful for higher frequency monitoring

`inverse` logical whether the barometric relationship is inverse

`role` character - the name of the role

`...` additional arguments

**Value**

barometric efficiency using Clark's method

**References**

Clark, W.E., 1967. Computing the barometric efficiency of a well. *Journal of the Hydraulics Division*, 93(4), pp.93-98.

Batu, V., 1998. *Aquifer hydraulics: a comprehensive guide to hydrogeologic data analysis*. John Wiley & Sons.

**See Also**

Other barometric: [step\\_baro\\_harmonic\(\)](#), [step\\_baro\\_least\\_squares\(\)](#)

**Examples**

```
data(kennel_2020)

clarks <- recipe(wl~., kennel_2020) |>
  step_baro_clark(wl, baro, lag_space = 1) |> # 1 minutes (every minute differences)
  step_baro_clark(wl, baro, lag_space = 60) |> # 60 minutes (hourly differences)
  step_baro_clark(wl, baro, lag_space = 1440) |> # 1440 minutes (daily differences)
  prep() |>
  bake()

clarks$get_step_data("barometric_efficiency")
```

---

```
step_baro_frequency_semi_confined
      step_baro_frequency_semi_confined
```

---

## Description

Work in Progress: Do not use. Rojstaczer 1988 solution  
 step\_baro\_frequency\_semi\_confined

## Usage

```
step_baro_frequency_semi_confined(
  .rec,
  frequency,
  radius_well,
  transmissivity,
  storage_confining,
  storage_aquifer,
  diffusivity_confining,
  diffusivity_vadose,
  thickness_confining,
  thickness_vadose,
  loading_efficiency,
  attenuation,
  role = "predictor",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>frequency</code>	the frequency of the response
<code>radius_well</code>	well radius
<code>transmissivity</code>	aquifer transmissivity (L*L/t)
<code>storage_confining</code>	confining layer storativity (L/L)
<code>storage_aquifer</code>	aquifer storativity (L/L)
<code>diffusivity_confining</code>	confining layer diffusivity
<code>diffusivity_vadose</code>	air diffusivity of vadose zone
<code>thickness_confining</code>	confining layer thickness

```

thickness_vadose      vadose thickness
loading_efficiency    the loading efficiency of the aquifer
attenuation           an attenuation factor
role                  character - the name of the role
...                   additional arguments

```

**Value**

complex response vector in frequency domain

---

```

step_baro_frequency_unconfined
      step_baro_frequency_semi_confined

```

---

**Description**

Work in Progress: Do not use. Rojstaczer 1988 solution  
*step\_baro\_frequency\_unconfined*

**Usage**

```

step_baro_frequency_unconfined(
  .rec,
  frequency,
  radius_well,
  storage_aquifer,
  specific_yield,
  k_vertical,
  diffusivity_vertical,
  diffusivity_vadose,
  thickness_saturated_well,
  thickness_vadose,
  thickness_aquifer,
  loading_efficiency,
  attenuation,
  role = "predictor",
  ...
)

```

**Arguments**

```

.rec      the R6 recipe object.
frequency the frequency of the response
radius_well well radius

```

**storage\_aquifer** aquifer storativity (L/L)  
**specific\_yield** the specific yeild of of the unconfined system  
**k\_vertical** vertical hydraulic conductivity of unconfined aquifer  
**diffusivity\_vertical** vertical diffusivity of unconfined aquifer  
**diffusivity\_vadose** air diffusivity of vadose zone  
**thickness\_saturated\_well** length of saturated well  
**thickness\_vadose** vadose thickness  
**thickness\_aquifer** aquifer thickness  
**loading\_efficiency** the loading efficiency of the aquifer  
**attenuation** an attenuation factor  
**role** character - the name of the role  
**...** additional arguments

**Value**

complex response vector in frequency domain

---

**step\_baro\_harmonic**    *step\_baro\_harmonic*

---

**Description**

Estimate the static barometric efficiency using harmonic methods (Ratio, Acworth, Rau, Transfer)

**Usage**

```

step_baro_harmonic(
  .rec,
  time,
  water_level,
  barometric_pressure,
  earth_tide,
  frequency = c(1.9324, 2),
  cycle_size = 86400,
  starting_value = 0,
  inverse = FALSE,
  role = "augment",
  ...
)
  
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	name of column that holds the time information
<code>water_level</code>	numeric vector of the dependent variable (ie:water level)
<code>barometric_pressure</code>	numeric vector of the independent variable (ie:barometric pressure)
<code>earth_tide</code>	variable unquoted Earth tide column name
<code>frequency</code>	numeric vector - the frequencies of the sin and cos curves
<code>cycle_size</code>	numeric - the period of the sin and cos curves
<code>starting_value</code>	numeric - the starting position of the sin and cos curves. This may be specified to have more control over the signal phase.
<code>inverse</code>	logical whether the barometric relationship is inverse
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

double barometric efficiency using different methods

**References**

Acworth, R.I., Halloran, L.J., Rau, G.C., Cuthbert, M.O. and Bernardi, T.L., 2016. An objective frequency domain method for quantifying confined aquifer compressible storage using Earth and atmospheric tides. *Geophysical Research Letters*, 43(22), pp.11-671.

**See Also**

Other barometric: [step\\_baro\\_clark\(\)](#), [step\\_baro\\_least\\_squares\(\)](#)

**Examples**

```
data("kennel_2020")
library(data.table)
library(collapse)

formula <- as.formula(wl~.)
frec = recipe(formula = formula, data = kennel_2020) |>
  step_baro_harmonic(datetime,
                     wl,
                     baro,
                     et,
                     inverse = FALSE)
```

---

```
step_baro_least_squares  
  step_baro_least_squares
```

---

## Description

Least squares solution for calculating barometric efficiency

## Usage

```
step_baro_least_squares(  
  .rec,  
  water_level,  
  barometric_pressure,  
  lag_space = 1L,  
  inverse = FALSE,  
  differences = FALSE,  
  role = "augment",  
  ...  
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>water_level</code>	numeric vector of the dependent variable (ie:water level)
<code>barometric_pressure</code>	numeric vector of the independent variable (ie:barometric pressure)
<code>lag_space</code>	integer spacing for lags, useful for higher frequency monitoring
<code>inverse</code>	logical whether the barometric relationship is inverse
<code>differences</code>	numeric vector number of samples between differences
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

barometric efficiency using least squares

## See Also

Other barometric: [step\\_baro\\_clark\(\)](#), [step\\_baro\\_harmonic\(\)](#)

## Examples

```
data(kennel_2020)

least_squares <- recipe(wl~., kennel_2020) |>
  step_baro_least_squares(wl, baro) |> # 1 minutes (every minute differences)
  step_baro_least_squares(wl, baro, lag_space = 1440, differences = TRUE) |>
  prep() |>
  bake()

least_squares$get_step_data("barometric_efficiency")
```

---

step_center	<i>step_center</i>
-------------	--------------------

---

## Description

Adds a step to center a data column(s)

## Usage

```
step_center(
  .rec,
  terms,
  role = "predictor",
  skip = FALSE,
  na_rm = TRUE,
  fun = collapse::fmean,
  keep_original_cols = FALSE,
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>na_rm</code>	logical - should NA values be removed from calculations
<code>fun</code>	function - the function that is applied to a list or columns of a <code>data.frame</code> like object.
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_center(x) |>
  prep() |>
  bake()
```

---

step_check_na	<i>step_check_na</i>
---------------	----------------------

---

**Description**

Check columns for NA

**Usage**

```
step_check_na(.rec, terms, role = "check", ...)
```

**Arguments**

<b>.rec</b>	the R6 recipe object.
<b>terms</b>	the unquoted names of the variables to use or a selector function. terms replaces the ‘...’ of the recipes package but requires variables to be included within ‘c()’. For example to include variables x and y you would write ‘c(x,y)’ in the hydrorecipes package.
<b>role</b>	character - the name of the role
<b>...</b>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_check_na(x) |>
  prep() |>
  bake()
```

---

```
step_check_spacing  step_check_spacing
```

---

### Description

Check the spacing of a variable

### Usage

```
step_check_spacing(.rec, terms, role = "check", ...)
```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

### Value

an updated recipe

### Examples

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_check_spacing(x) |>
  prep() |>
  bake()
```

---

```
step_compare_columns  step_check_spacing
```

---

### Description

Check the spacing of a variable

**Usage**

```
step_compare_columns(  
  .rec,  
  data,  
  compare,  
  role = "add",  
  n_sd = 4,  
  na_rm = TRUE,  
  ...  
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>data</code>	variable unquoted data column name
<code>compare</code>	variable unquoted column name for comparison
<code>role</code>	character - the name of the role
<code>n_sd</code>	numeric - number of standard deviations for the scaling
<code>na_rm</code>	logical - should NA values be removed from calculations
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
data("kennel_2020")  
  
kennel_2020[1e4, w1 := 13.36]  
  
frec = recipe(w1~baro, data = kennel_2020) |>  
  step_compare_columns(data = w1, compare = baro, n_sd = 15) |>  
  prep() |>  
  bake()
```

---

`step_convolve_exponential`

*step\_convolve\_exponential*

---

**Description**

linearly convolve a gamma kernel with a data series.

**Usage**

```
step_convolve_exponential(
  .rec,
  terms,
  amplitude,
  theta,
  align = "right",
  max_length = Inf,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>amplitude</code>	amplitude
<code>theta</code>	scale
<code>align</code>	character center, left or right align the convolution
<code>max_length</code>	the maximum length of the kernel
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
formula <- as.formula(x~y+z)
rows <- 1e4

dat <- data.frame(x = rep(1, rows),
                 y = 1:rows,
                 z = cumsum(rnorm(rows)))

frec = recipe(formula = formula, data = dat) |>
  step_convolve_gamma(z, amplitude = 1, theta = 1, k = 1) |>
  plate("tbl")
```

---

```
step_convolve_gamma  step_convolve_gamma
```

---

## Description

linearly convolve a gamma kernel with a data series.

## Usage

```
step_convolve_gamma(
  .rec,
  terms,
  amplitude,
  k,
  theta,
  align = "right",
  max_length = Inf,
  role = "predictor",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>amplitude</code>	amplitude
<code>k</code>	shape
<code>theta</code>	scale
<code>align</code>	character center, left or right align the convolution
<code>max_length</code>	the maximum length of the kernel
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

**Examples**

```

formula <- as.formula(x~y+z)
rows <- 1e4

dat <- data.frame(x = rep(1, rows),
                  y = 1:rows,
                  z = cumsum(rnorm(rows)))

frec = recipe(formula = formula, data = dat) |>
  step_convolve_gamma(z, amplitude = 1, theta = 1, k = 1) |>
  plate("tbl")

```

---

```

step_cross_correlation
      step_cross_correlation

```

---

**Description**

Calculate the autocorrelation function or cross-correlation

**Usage**

```
step_cross_correlation(.rec, terms, lag_max = 100, role = "predictor", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipies package.
<code>lag_max</code>	maximum lag to calculate
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```

formula <- as.formula(y~x)
rows <- 1e4

dat <- data.frame(x = rnorm(rows),
                  y = as.numeric(1:rows))

frec = recipe(formula = formula, data = dat) |>
  step_cross_correlation(x)

frec = recipe(formula = formula, data = dat) |>
  step_cross_correlation(c(x, y))

```

---

step\_distributed\_lag *step\_distributed\_lag*

---

**Description**

Generates distributed lag vectors. For regular spaced lags this uses an FFT based method which is faster and more memory efficient.

**Usage**

```

step_distributed_lag(
  .rec,
  terms,
  n_lag = 12L,
  lag_max = 86400L,
  knots = NA_real_,
  basis_matrix = NA_real_,
  intercept = FALSE,
  role = "predictor",
  ...
)

```

**Arguments**

<b>.rec</b>	the R6 recipe object.
<b>terms</b>	the unquoted names of the variables to use or a selector function. terms replaces the ‘...’ of the recipes package but requires variables to be included within ‘c()’. For example to include variables x and y you would write ‘c(x,y)’ in the hydrorecipes package.
<b>n_lag</b>	number of lags to calculate (enter either lag_max and n_lag or knots)
<b>lag_max</b>	maximum lag to calculate (enter either lag_max and n_lag or knots)
<b>knots</b>	specify the knot locations

<code>basis_matrix</code>	user specified <code>basis_matrix</code>
<code>intercept</code>	include intercept in basis matrix
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**References**

Gasparri, A., 2011. Distributed Lag Linear and Non-Linear Models in R: The Package `dlm`. *Journal of statistical software* 465 43, 1–20.

**Examples**

```
formula <- as.formula(y~x)
rows <- 1e4

dat <- data.frame(x = rnorm(rows),
                 y = as.numeric(1:rows),
                 z = rnorm(rows))

frec = recipe(formula = formula, data = dat) |>
  step_distributed_lag(x, knots = hydrorecipes::log_lags_arma(6, 800))
```

---

<code>step_drop_columns</code>	<i>step_drop_columns</i>
--------------------------------	--------------------------

---

**Description**

Removes regressors from recipe.

**Usage**

```
step_drop_columns(.rec, terms, role = "modify", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
rows <- 20
formula <- as.formula(y~x)

dat <- data.frame(x = rnorm(rows),
                  y = as.numeric(1:rows),
                  z = rnorm(rows))

frec = recipe(formula = formula, data = dat) |>
  step_drop_columns(x)
```

---

 step\_dummy

*step\_dummy*


---

**Description**

dummy encode a factor or factor like variable.

**Usage**

```
step_dummy(
  .rec,
  terms,
  one_hot = FALSE,
  role = "predictor",
  skip = FALSE,
  keep_original_cols = FALSE,
  ...
)
```

**Arguments**

<b>.rec</b>	the R6 recipe object.
<b>terms</b>	the unquoted names of the variables to use or a selector function. terms replaces the ‘...’ of the recipes package but requires variables to be included within ‘c()’. For example to include variables x and y you would write ‘c(x,y)’ in the hydrorecipes package.
<b>one_hot</b>	logical - use one hot encoding.
<b>role</b>	character - the name of the role
<b>skip</b>	logical - should the step be skipped
<b>keep_original_cols</b>	logical - keep the original columns or replace them
<b>...</b>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = factor(sample(1:10, 100, replace = TRUE)),
                 y = rnorm(100))

rec <- recipe(y~x, data = dat) |>
  step_dummy(x, one_hot = FALSE)
rec <- recipe(y~x, data = dat) |>
  step_dummy(x, one_hot = TRUE)
```

---

step_earthtide	<i>step_earthtide</i>
----------------	-----------------------

---

**Description**

Generate synthetic Earth tide waves and wave groups.

**Usage**

```
step_earthtide(
  .rec,
  terms,
  do_predict = TRUE,
  method = "gravity",
  latitude = 0,
  longitude = 0,
  elevation = 0,
  azimuth = 0,
  gravity = 0,
  earth_radius = 6378136.3,
  earth_eccen = 0.0066943979514,
  cutoff = 1e-06,
  catalog = "ksm04",
  eop = NULL,
  scale = TRUE,
  n_thread = 1L,
  astro_update = 1L,
  interp_factor = 1L,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipes package.
<code>do_predict</code>	run in predict or analyze mode
<code>method</code>	One or more of "gravity", "tidal_potential", "tidal_tilt", "vertical_displacement", "horizontal_displacement", "n_s_displacement", "e_w_displacement", "vertical_strain", "areal_strain", "volume_strain", "horizontal_strain", or "ocean_tides", "pole_tide", "lod_tide". The pole tide and lod_tide are used in predict mode even if <code>do_predict</code> is FALSE. More than one value can only be used if <code>do_predict == TRUE</code> .
<code>latitude</code>	The station latitude (numeric) defaults to 0.
<code>longitude</code>	The station longitude (numeric) defaults to 0.
<code>elevation</code>	The station elevation (m) (numeric) defaults to 0.
<code>azimuth</code>	Earth azimuth (numeric) defaults to 0.
<code>gravity</code>	Gravity at the station ( $m/s^2$ ) (numeric) 0 to estimate gravity from elevation and latitude.
<code>earth_radius</code>	Radius of earth (m) (numeric) defaults to 6378136.3
<code>earth_eccen</code>	Eccentricity of earth (numeric) defaults to 6.69439795140e-3
<code>cutoff</code>	Cutoff amplitude for constituents (numeric) defaults to 1e-6.
<code>catalog</code>	Use the "hw95s" catalog or "ksm04" catalog (character).
<code>eop</code>	User defined Earth Orientation Parameter (EOP) data.frame with the following columns: <code>datetime</code> , <code>ddt</code> , <code>ut1_utc</code> , <code>lod</code> , <code>x</code> , <code>y</code> , <code>dx</code> , <code>dy</code>
<code>scale</code>	Scale results when <code>do_predict</code> is FALSE
<code>n_thread</code>	Number of threads to use for parallel processing (integer).
<code>astro_update</code>	How often to update astro parameters in number of samples. This speeds up code but may make it slightly less accurate.
<code>interp_factor</code>	calculate for every <code>interp_factor</code> samples. This may be faster while being more accurate than adjusting the cutoff.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```

library(data.table)

data(kennel_2020)
latitude   <- 34.23411           # latitude
longitude  <- -118.678          # longitude
elevation  <- 500                # elevation
cutoff     <- 1e-5              # cutoff
catalog    <- 'ksm04'           # hartmann wenzel catalog
astro_update <- 300              # how often to update astro parameters
method     <- 'volume_strain'   # which potential to calculate

wave_groups_dl <- data.table::as.data.table(earthtide::eterna_wavegroups)
wave_groups_dl <- na.omit(wave_groups_dl[time == "1 month"])
wave_groups_dl <- wave_groups_dl[wave_groups_dl$start > 0.5,]
wave_groups_dl <- wave_groups_dl[, list(start, end)]
ngr <- nrow(wave_groups_dl)

rec <- recipe(wl~baro+datetime, data = kennel_2020) |>
  step_earthtide(datetime,
    wave_groups = wave_groups_dl,
    latitude = latitude,
    longitude = longitude,
    elevation = elevation,
    cutoff = cutoff,
    catalog = catalog)

```

---

```
step_fft_coherence  step_fft_coherence
```

---

**Description**

estimates the coherence between terms.

**Usage**

```
step_fft_coherence(.rec, terms, role = "augment", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipes package.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
data(kennel_2020)

form <- as.formula("wl~.")

formula <- as.formula(wl~baro + et)
frec = recipe(formula = formula, data = kennel_2020) |>
  step_fft_pgram(c(wl, baro, et), spans = 7) |>
  step_fft_coherence() |>
  prep() |>
  bake()
```

---

step_fft_pgram	<i>step_fft_pgram</i>
----------------	-----------------------

---

**Description**

Periodgrams and cross-periodograms using a method similar to `stats::spec.pgram`.

**Usage**

```
step_fft_pgram(
  .rec,
  terms,
  spans = 3,
  detrend = TRUE,
  demean = TRUE,
  lst = TRUE,
  taper = 0.1,
  pad_fft = TRUE,
  time_step = 1,
  role = "predictor",
  ...
)
```

**Arguments**

<b>.rec</b>	the R6 recipe object.
<b>terms</b>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.

<code>spans</code>	vector of odd integers giving the widths of modified Daniell smoothers to be used to smooth the periodogram.
<code>detrend</code>	logical. If <code>TRUE</code> , remove a linear trend from the series. This will also remove the mean.
<code>demean</code>	logical. If <code>TRUE</code> , subtract the mean of the series.
<code>lst</code>	logical return a list?
<code>taper</code>	specifies the proportion of data to taper. A split cosine bell taper is applied to this proportion of the data at the beginning and end of the series.
<code>pad_fft</code>	logical Zero pad the list for faster FFT calculation?
<code>time_step</code>	numeric monitoring interval size
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
formula <- as.formula(y~.)

dat <- data.frame(x = rnorm(200),
                 y = rnorm(200))

frec = recipe(formula = formula, data = dat) |>
  step_fft_pgram(c(x,y))
```

---

```
step_fft_transfer_experimental
  step_fft_transfer_experimental
```

---

**Description**

Calculates the transfer function with pgram results.

**Usage**

```
step_fft_transfer_experimental(
  .rec,
  terms,
  spans = 3,
  detrend = TRUE,
  demean = TRUE,
  taper = 0.1,
```

```

    n_groups = 200,
    time_step = 1,
    formula = NULL,
    role = "augment",
    ...
  )

```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>spans</code>	vector of odd integers giving the widths of modified Daniell smoothers to be used to smooth the periodogram.
<code>detrend</code>	logical. If <code>TRUE</code> , remove a linear trend from the series. This will also remove the mean.
<code>demean</code>	logical. If <code>TRUE</code> , subtract the mean of the series.
<code>taper</code>	specifies the proportion of data to taper. A split cosine bell taper is applied to this proportion of the data at the beginning and end of the series.
<code>n_groups</code>	number of results
<code>time_step</code>	numeric monitoring interval size
<code>formula</code>	formula notation to specify inputs and outputs
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

## Examples

```

data(kennel_2020)

form <- as.formula("wl~.")

rec <- recipe(form, kennel_2020) |>
  step_fft_transfer_experimental(c(wl, baro, et), spans = 3) |>
  plate()

```

---

```
step_fft_transfer_pgram
  step_fft_transfer_pgram
```

---

## Description

Calculates the transfer function with pgram results.

## Usage

```
step_fft_transfer_pgram(
  .rec,
  terms,
  spans = 3,
  detrend = TRUE,
  demean = TRUE,
  taper = 0.1,
  time_step = 1,
  formula = NULL,
  role = "augment",
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>spans</code>	vector of odd integers giving the widths of modified Daniell smoothers to be used to smooth the periodogram.
<code>detrend</code>	logical. If <code>TRUE</code> , remove a linear trend from the series. This will also remove the mean.
<code>demean</code>	logical. If <code>TRUE</code> , subtract the mean of the series.
<code>taper</code>	specifies the proportion of data to taper. A split cosine bell taper is applied to this proportion of the data at the beginning and end of the series.
<code>time_step</code>	numeric monitoring interval size
<code>formula</code>	formula notation to specify inputs and outputs
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

**Examples**

```
data(kennel_2020)

form <- as.formula("wl~.")

rec <- recipe(form, kennel_2020) |>
  step_fft_transfer_pgram(c(wl, baro, et), spans = 3) |>
  plate()
```

---

```
step_fft_transfer_welch
  step_fft_transfer_welch
```

---

**Description**

calculates the transfer function with Welch's results.

**Usage**

```
step_fft_transfer_welch(
  .rec,
  terms,
  length_subset,
  overlap = 0.8,
  window,
  time_step = 1,
  formula = NULL,
  role = "augment",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>length_subset</code>	length of fft section
<code>overlap</code>	amount of overlap
<code>window</code>	window weights
<code>time_step</code>	numeric monitoring interval size
<code>formula</code>	formula notation to specify inputs and outputs
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
data(kennel_2020)
form <- as.formula("wl~.")

rec <- recipe(form, kennel_2020) |>
step_fft_transfer_welch(c(wl, baro, et),
                        length_subset = 1440*8 + 1,
                        overlap = 0.6,
                        window = window_nuttall(1440*8+1)) |>
plate()
```

---

<code>step_fft_welch</code>	<i>step_fft_welch</i>
-----------------------------	-----------------------

---

**Description**

calculates the periodogram (estimate of spectral density) using Welch's method.

**Usage**

```
step_fft_welch(
  .rec,
  terms,
  length_subset,
  overlap = 0.8,
  window,
  time_step = 1,
  role = "augment",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipes package.
<code>length_subset</code>	length of fft section
<code>overlap</code>	amount of overlap
<code>window</code>	window weights

<code>time_step</code>	numeric monitoring interval size
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
formula <- as.formula(y~.)

dat <- data.frame(x = rnorm(200),
                 y = rnorm(200))

frec = recipe(formula = formula, data = dat) |>
  step_fft_welch(c(x,y), length_subset = 10, window = window_rectangle(10))
```

---

<code>step_find_interval</code>	<i>step_find_interval</i>
---------------------------------	---------------------------

---

**Description**

divides a series into intervals and then performs dummy encoding.

**Usage**

```
step_find_interval(.rec, terms, vec, role = "augment", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipes package.
<code>vec</code>	a vector of break points
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

## Examples

```
formula <- as.formula(y~x)

rows = 200
dat <- data.frame(x = rnorm(rows),
                  y = 1:rows,
                  z = rnorm(rows))

frec1 = recipe(formula = formula, data = dat) |>
  step_find_interval(x, vec = c(-0.1, 0.0, 0.1)) |>
  plate("tbl")
```

---

step_harmonic	<i>step_harmonic</i>
---------------	----------------------

---

## Description

Add sin and cos terms for harmonic analysis

## Usage

```
step_harmonic(
  .rec,
  terms,
  frequency = NA_real_,
  cycle_size = NA_real_,
  starting_value = 0,
  role = "predictor",
  skip = FALSE,
  keep_original_cols = FALSE,
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipes package.
<code>frequency</code>	numeric vector - the frequencies of the sin and cos curves
<code>cycle_size</code>	numeric - the period of the sin and cos curves
<code>starting_value</code>	numeric - the starting position of the sin and cos curves. This may be specified to have more control over the signal phase.

<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = 1:10, y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_harmonic(x,
               frequency = 2.0,
               cycle_size = 4.0,
               starting_value = 0.0)
```

---

<code>step_intercept</code>	<i>step_intercept</i>
-----------------------------	-----------------------

---

**Description**

Add an intercept term

**Usage**

```
step_intercept(.rec, terms, value = 1, role = "predictor", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipies</code> package.
<code>value</code>	what value to use (typically 1.0)
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = 1:10, y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_intercept()
```

---

```
step_kernel_divide_naive
  step_kernel_divide_naive
```

---

**Description**

Divide a signal by a kernel in the frequency domain.

**Usage**

```
step_kernel_divide_naive(.rec, terms, kernel, role = "predictor", ...)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>kernel</code>	the convolution kernel
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

---

```
step_kernel_filter  step_kernel_filter
```

---

**Description**

linearly convolve a kernel with a data series.

## Usage

```
step_kernel_filter(  
  .rec,  
  terms,  
  kernel,  
  align = "center",  
  role = "predictor",  
  ...  
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>kernel</code>	the convolution kernel
<code>align</code>	character center, left or right align the convolution
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

## Examples

```
formula <- as.formula(x~y+z)  
rows <- 1e4  
  
dat <- data.frame(x = rep(1, rows),  
                 y = 1:rows,  
                 z = cumsum(rnorm(rows)))  
  
frec = recipe(formula = formula, data = dat) |>  
  step_kernel_filter(z, kernel = list(rep(1, 1001)/1001), align = "center") |>  
  plate("tbl")
```

---

```
step_lead_lag      step_lead_lag
```

---

## Description

Lag or lead a column or columns. This requires a sorted and regular time series.

## Usage

```
step_lead_lag(
  .rec,
  terms,
  lag,
  n_shift = 0L,
  n_subset = 1L,
  role = "predictor",
  skip = FALSE,
  keep_original_cols = FALSE,
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>lag</code>	integer vector - number of samples to lag or lead. Negative numbers indicate leading a vector.
<code>n_shift</code>	integer - number of values to shift the starting position when <code>n_subset</code> is not equal to 0. The value of <code>n_shift</code> has to be less than <code>'n_subset'</code> .
<code>n_subset</code>	integer - spacing between adjacent samples in the result.
<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

## Value

an updated recipe

**Examples**

```

dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_lead_lag(x, lag = 1)

rec <- recipe(y~x, data = dat) |>
  step_lead_lag(x, lag = 1, n_subset = 5)

rec <- recipe(y~x, data = dat) |>
  step_lead_lag(x, lag = 1, n_shift = 2, n_subset = 5)

```

---

step_multiply	<i>step_multiply</i>
---------------	----------------------

---

**Description**

step\_multiply

**Usage**

```

step_multiply(
  .rec,
  terms,
  values = 1,
  role = "predictor",
  skip = FALSE,
  keep_original_cols = FALSE,
  ...
)

```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipies</code> package.
<code>values</code>	multiply column(s) by these values
<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_multiply(x, value = 4)
```

---

step\_nls

*step\_nls*

---

**Description**

Uses the Eigen C++ library fast versions to generate predictions and coefficients from a recipe.

**Usage**

```
step_nls(
  .rec,
  formula,
  algorithm = "lm",
  n_subset = 1L,
  n_shift = 0L,
  range = c(-Inf, Inf),
  control = gsl_nls_control(xtol = 1e-08),
  trace = FALSE,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>formula</code>	formula for the regression
<code>algorithm</code>	character string specifying the algorithm to use. The following choices are supported: <ul style="list-style-type: none"> <li>• "lm" Levenberg-Marquardt algorithm (default).</li> <li>• "lmaccel" Levenberg-Marquardt algorithm with geodesic acceleration. Stability is controlled by the <code>avmax</code> parameter in <code>control</code>, setting <code>avmax</code> to zero is analogous to not using geodesic acceleration.</li> <li>• "dogleg" Powell's dogleg algorithm.</li> </ul>

- "ddogleg" Double dogleg algorithm, an improvement over "dogleg" by including information about the Gauss-Newton step while the iteration is still far from the minimum.
- "subspace2D" 2D generalization of the dogleg algorithm. This method searches a larger subspace for a solution, it can converge more quickly than "dogleg" on some problems.

n_subset	integer - spacing between adjacent samples in the result.
n_shift	integer - number of values to shift the starting position when n_subset is not equal to 0. The value of n_shift has to be less than 'n_subset'.
range	limit the fitting range to observations between range[1] and range[2]
control	an optional list of control parameters to tune the least squares iterations and multistart algorithm. See <a href="#">gsl_nls_control</a> for the available control parameters and their default values.
trace	logical value indicating if a trace of the iteration progress should be printed. Default is FALSE. If TRUE, the residual (weighted) sum-of-squares and the current parameter estimates are printed after each iteration.
role	character - the name of the role
...	additional arguments

**Value**

an updated recipe

**See Also**

Other ols: [step\\_ols\(\)](#)

**Examples**

```
data("kennel_2020")
kennel_2020[, datetime := as.numeric(datetime)]
formula <- as.formula(wl~.)
n_knots <- 12
deg_free <- 27
max_lag <- 1 + 720

frec = recipe(formula = formula, data = unclass(kennel_2020)) |>
  step_distributed_lag(baro, knots = hydrorecipes::log_lags_arma(n_knots, max_lag)) |>
  step_spline_b(datetime, df = deg_free, intercept = FALSE) |>
  step_intercept() |>
  step_drop_columns(baro) |>
  step_drop_columns(datetime) |>
  step_ols(formula) |>
  prep() |>
  bake()
```

---

step_normalize	<i>step_normalize</i>
----------------	-----------------------

---

## Description

step\_normalize

## Usage

```
step_normalize(  
  .rec,  
  terms,  
  role = "predictor",  
  skip = FALSE,  
  na_rm = TRUE,  
  keep_original_cols = FALSE,  
  ...  
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipies</code> package.
<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>na_rm</code>	logical - should NA values be removed from calculations
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

## Value

an updated recipe

## Examples

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))  
  
rec <- recipe(y~x, data = dat) |>  
  step_normalize(x)
```

---

step_ols	<i>step_ols</i>
----------	-----------------

---

## Description

Uses the Eigen C++ library fast versions to generate predictions and coefficients from a recipe.

## Usage

```
step_ols(.rec, formula, role = "predictor", do_response = TRUE, ...)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>formula</code>	formula for the regression
<code>role</code>	character - the name of the role
<code>do_response</code>	logical calculate and return the responses?
<code>...</code>	additional arguments

## Value

an updated recipe

## See Also

Other ols: [step\\_nls\(\)](#)

## Examples

```
data("kennel_2020")
kennel_2020[, datetime := as.numeric(datetime)]
formula <- as.formula(wl~.)
n_knots <- 12
deg_free <- 27
max_lag <- 1 + 720

frec = recipe(formula = formula, data = unclass(kennel_2020)) |>
  step_distributed_lag(baro, knots = hydrorecipes::log_lags_arma(n_knots, max_lag)) |>
  step_spline_b(datetime, df = deg_free, intercept = FALSE) |>
  step_intercept() |>
  step_drop_columns(baro) |>
  step_drop_columns(datetime) |>
  step_ols(formula) |>
  prep() |>
  bake()
```

---

```
step_ols_gap_fill      step_ols_gap_fill
```

---

### Description

```
step_ols_gap_fill
```

### Usage

```
step_ols_gap_fill(.rec, terms, recipe, role = "predictor", ...)
```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>recipe</code>	Recipe to use for filling gaps
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

### Value

an updated recipe

### Examples

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))
```

---

```
step_pca              step_pca
```

---

### Description

'StepPca' Does PCA for a set of columns. This currently is an in house function. Use at your own risk!

**Usage**

```
step_pca(
  .rec,
  terms,
  na_rm = TRUE,
  n_comp = 3,
  center = TRUE,
  scale = TRUE,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipies</code> package.
<code>na_rm</code>	logical - should NA values be removed from calculations
<code>n_comp</code>	number of components to retain
<code>center</code>	center values before PCA
<code>scale</code>	scale values before PCA
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**Examples**

```
set.seed(1)

formula <- as.formula(x~a+b+d+e+f+g)
rows <- 1000

dat <- data.frame(x = rnorm(rows),
                 a = rnorm(rows),
                 b = rnorm(rows),
                 d = rnorm(rows),
                 e = rnorm(rows),
                 f = rnorm(rows),
                 g = rnorm(rows))

rec = recipe(formula = formula, data = dat) |>
  step_pca(c(x,a,b,d,e,f,g)) |>
  plate()
```

---

step_scale	<i>step_scale</i>
------------	-------------------

---

## Description

Adds a step to scale a data column(s)

## Usage

```
step_scale(
  .rec,
  terms,
  role = "predictor",
  skip = FALSE,
  na_rm = TRUE,
  fun = collapse::fsd,
  n_sd = 1L,
  keep_original_cols = FALSE,
  ...
)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>skip</code>	logical - should the step be skipped
<code>na_rm</code>	logical - should NA values be removed from calculations
<code>fun</code>	function - the function that is applied to a list or columns of a <code>data.frame</code> like object.
<code>n_sd</code>	numeric - number of standard deviations for the scaling
<code>keep_original_cols</code>	logical - keep the original columns or replace them
<code>...</code>	additional arguments

## Value

an updated recipe

**Examples**

```
dat <- data.frame(x = rnorm(10), y = rnorm(10))

rec <- recipe(y~x, data = dat) |>
  step_scale(x)
```

---

```
step_slug_cbp      step_slug_cbp
```

---

**Description**

Cooper, Bredehoeft and Papadopoulos, 1967 Slug test solution

**Usage**

```
step_slug_cbp(
  .rec,
  time,
  radius = 1,
  radius_casing = 0.15,
  radius_well = 0.15,
  specific_storage = 1e-06,
  hydraulic_conductivity = 1e-04,
  head_0 = 1,
  thickness = 1,
  n_terms = 16,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	the time for evaluation (t)
<code>radius</code>	distance from line source or center of well
<code>radius_casing</code>	radius of casing
<code>radius_well</code>	radius of well screen
<code>specific_storage</code>	specific storage of aquifer (L/L)
<code>hydraulic_conductivity</code>	the hydraulic conductivity (L/t)
<code>head_0</code>	initial displacement
<code>thickness</code>	the aquifer thickness (L)
<code>n_terms</code>	number of terms for laplace solution inversion
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**References**

Cooper, H.H., J.D. Bredehoeft and S.S. Papadopoulos, 1967. Response of a finite-diameter well to an instantaneous charge of water, Water Resources Research, vol. 3, no. 1, pp. 263-269.

**Examples**

```
# check vs. CBP 1967 table 1
times <- rep(c(1.0, 2.15, 4.64), 4) * 10^(rep(c(-3, -2, -1, 0), each = 3))
dat <- list(x = times)

formula = formula(x~.)

frec1 = recipe(formula = formula, data = dat) |>
  step_slug_cbp(
    time = x,
    radius = 1.0,
    radius_casing = 1.0,
    radius_well = 1.0,
    specific_storage = 1e-1,
    hydraulic_conductivity = 1.0,
    thickness = 1.0,
    head_0 = 1.0,
    n_terms = 12L
  ) |>
  plate("dt")
```

---

step\_spline\_b

*step\_spline\_b*

---

**Description**

generates basis splines.

**Usage**

```
step_spline_b(
  .rec,
  terms,
  df = 0L,
  internal_knots = NULL,
  boundary_knots = NULL,
  intercept = FALSE,
  periodic = FALSE,
  degree = 3L,
```

```

    role = "predictor",
    ...
  )

```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>df</code>	Degree of freedom that equals to the column number of the returned matrix. One can specify <code>df</code> rather than <code>knots</code> , then the function chooses <code>df - degree - as.integer(intercept)</code> internal knots at suitable quantiles of <code>x</code> ignoring missing values and those <code>x</code> outside of the boundary. For periodic splines, <code>df - as.integer(intercept)</code> internal knots will be chosen at suitable quantiles of <code>x</code> relative to the beginning of the cyclic intervals they belong to (see Examples) and the number of internal knots must be greater or equal to the specified <code>degree - 1</code> . If internal knots are specified via <code>knots</code> , the specified <code>df</code> will be ignored.
<code>internal_knots</code>	equivalent to <code>knots</code> from <code>'splines2::bSplines'</code>
<code>boundary_knots</code>	equivalent to <code>Boundary.knots</code> from <code>'splines2::bSplines'</code>
<code>intercept</code>	If <code>TRUE</code> , the complete basis matrix will be returned. Otherwise, the first basis will be excluded from the output.
<code>periodic</code>	A logical value. If <code>TRUE</code> , the periodic splines will be returned. The default value is <code>FALSE</code> .
<code>degree</code>	A nonnegative integer specifying the degree of the piecewise polynomial. The default value is 3 for cubic splines. Zero degree is allowed for piecewise constant basis functions.
<code>role</code>	character - the name of the role
<code>...</code>	Optional arguments that are not used.

### Value

an updated recipe

### Examples

```

formula <- as.formula(x-y+z)
rows <- 1e5

dat <- data.frame(x = rnorm(rows),
                 y = 1:rows,
                 z = cumsum(rnorm(rows)))
ik <- collapse::fquantile(dat$x, probs = seq(0, 1, 0.1))
bk <- ik[c(1, length(ik))]

```

```

ik <- ik[-c(1, length(ik))]

frec = recipe(formula = formula, data = dat) |>
  step_spline_b(x, df = 11L, intercept = FALSE) |>
  plate("tbl")

```

---

```

step_spline_n      step_spline_n

```

---

## Description

generates basis splines.

## Usage

```

step_spline_n(
  .rec,
  terms,
  df = 0L,
  internal_knots = NULL,
  boundary_knots = NULL,
  intercept = FALSE,
  periodic = FALSE,
  degree = 3L,
  role = "predictor",
  ...
)

```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>df</code>	Degree of freedom that equals to the column number of returned matrix. One can specify <code>df</code> rather than <code>knots</code> , then the function chooses <code>df - 1 - as.integer(intercept)</code> internal knots at suitable quantiles of <code>x</code> ignoring missing values and those <code>x</code> outside of the boundary. Thus, <code>df</code> must be greater than or equal to 2. If internal knots are specified via <code>knots</code> , the specified <code>df</code> will be ignored.
<code>internal_knots</code>	equivalent to <code>knots</code> from <code>'splines2::bSplines'</code>
<code>boundary_knots</code>	equivalent to <code>Boundary.knots</code> from <code>'splines2::bSplines'</code>

<code>intercept</code>	If TRUE, the complete basis matrix will be returned. Otherwise, the first basis will be excluded from the output.
<code>periodic</code>	A logical value. If TRUE, the periodic splines will be returned. The default value is FALSE.
<code>degree</code>	A nonnegative integer specifying the degree of the piecewise polynomial. The default value is 3 for cubic splines. Zero degree is allowed for piecewise constant basis functions.
<code>role</code>	character - the name of the role
<code>...</code>	Optional arguments that are not used.

### Value

an updated recipe

### Examples

```
formula <- as.formula(x~y+z)
rows <- 1e5

dat <- data.frame(x = rnorm(rows),
                 y = 1:rows,
                 z = cumsum(rnorm(rows)))
ik <- collapse::fquantile(dat$x, probs = seq(0, 1, 0.1))
bk <- ik[c(1, length(ik))]
ik <- ik[-c(1, length(ik))]

frec = recipe(formula = formula, data = dat) |>
  step_spline_n(x, df = 11L, intercept = FALSE) |>
  plate("tbl")
```

---

`step_subset_na_omit` *step\_subset\_na\_omit*

---

### Description

selects rows from output.

### Usage

```
step_subset_na_omit(.rec, terms, role = "modify", ...)
```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the recipes package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the hydrorecipies package.

role character - the name of the role  
 ... additional arguments

### Value

an updated recipe

---

step\_subset\_rows *step\_subset\_rows*

---

### Description

selects rows from output.

### Usage

```
step_subset_rows(.rec, row_numbers, role = "modify", ...)
```

### Arguments

.rec the R6 recipe object.  
 row\_numbers integer vector of row numbers to keep.  
 role character - the name of the role  
 ... additional arguments

### Value

an updated recipe

### Examples

```
dat <- data.frame(x = as.numeric(1:200), y = rnorm(200))
formula <- as.formula(y~x)

frec1 = recipe(formula = formula, data = dat) |>
  step_subset_rows(row_numbers = c(1, 5, 10)) |>
  plate("dt")
```

---

```
step_subset_sample  step_subset_sample
```

---

### Description

selects rows from output.

### Usage

```
step_subset_sample(.rec, terms, size, role = "modify", ...)
```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'..'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipies</code> package.
<code>size</code>	number of samples.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

### Value

an updated recipe

---

```
step_transport_fractures_heat
      step_transport_fractures_heat
```

---

### Description

Sudicky and Frind 1982 solution adapted for heat. Two parallel fractures.

### Usage

```
step_transport_fractures_heat(
  .rec,
  time,
  distance_fracture,
  distance_matrix,
  temperature_influent = 15,
  time_influent = 0,
  temperature_initial = 10,
  fracture_aperture = 2e-04,
```

```

fracture_spacing = 1,
velocity = 0.1/86400,
thermal_conductivity_water = 0.615,
thermal_conductivity_solids = 3.4,
specific_heat_water = 4192,
specific_heat_solids = 908,
density_water = 1,
density_solids = 2.5,
porosity = 0.1,
n_terms = 30L,
role = "predictor",
...
)

```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>time</code>	vector elapsed time (t)
<code>distance_fracture</code>	vector distance along fracture (z)
<code>distance_matrix</code>	vector distance into matrix (x)
<code>temperature_influent</code>	vector temperature history (t_in)
<code>time_influent</code>	vector time of influent values (t_in)
<code>temperature_initial</code>	double temperature (t_0)
<code>fracture_aperture</code>	double fracture aperture (2b)
<code>fracture_spacing</code>	double fracture aperture (2B)
<code>velocity</code>	double water velocity in fracture (v)
<code>thermal_conductivity_water</code>	double water thermal conductivity ( _f)
<code>thermal_conductivity_solids</code>	double solids thermal conductivity ( _s)
<code>specific_heat_water</code>	double specific heat of water
<code>specific_heat_solids</code>	double specific heat of solid particles
<code>density_water</code>	double density of the water ( _w)
<code>density_solids</code>	double density of the solid particles ( _s)
<code>porosity</code>	double matrix porosity ( )
<code>n_terms</code>	integer the number of laplace terms
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**See Also**

Other transport: [step\\_transport\\_fractures\\_solute\(\)](#), [step\\_transport\\_ogata\\_banks\(\)](#)

**Examples**

```
formula <- as.formula(~time+z+x)

dat <- as.data.frame(expand.grid(10^(3:8),
                               seq(0.0, 100, 1),
                               c(0.0, 0.05)))

names(dat) <- c("time", "z", "x")

frec1 = recipe(formula = formula, data = dat) |>
  step_transport_fractures_heat(time = time,
                               distance_fracture = z,
                               distance_matrix = x) |>
  plate()
```

---

```
step_transport_fractures_solute
  step_transport_fractures_solute
```

---

**Description**

Sudicky and Frind 1982 solution. Two parallel fractures

**Usage**

```
step_transport_fractures_solute(
  .rec,
  time,
  distance_fracture,
  distance_matrix,
  concentration_influent = 1,
  time_influent = 0,
  concentration_initial = 0,
  fracture_aperture = 2e-04,
  fracture_spacing = 1,
  velocity = 0.1/86400,
  dispersivity_longitudinal = 0.1,
  diffusion = 1e-09,
```

```

sorption_fracture = 0,
sorption_matrix = 0,
decay = 1e+15,
density_bulk = 2.5,
porosity = 0.1,
tortuosity = 0.1,
n_terms = 30L,
role = "predictor",
...
)

```

### Arguments

<code>.rec</code>	the R6 recipe object.
<code>time</code>	vector elapsed time (t)
<code>distance_fracture</code>	vector distance along fracture (z)
<code>distance_matrix</code>	vector distance into matrix (x)
<code>concentration_influent</code>	vector concentration history (c_in)
<code>time_influent</code>	vector concentration history (t_in)
<code>concentration_initial</code>	double concentration (c_0)
<code>fracture_aperture</code>	double fracture aperture (2b)
<code>fracture_spacing</code>	double fracture aperture (2B)
<code>velocity</code>	double water velocity in fracture (v)
<code>dispersivity_longitudinal</code>	double longitudinal dispersivity ( _l)
<code>diffusion</code>	double free-water diffusion coefficient (D*)
<code>sorption_fracture</code>	double fracture distribution coefficient (K_f)
<code>sorption_matrix</code>	double matrix distribution coefficient (K_m)
<code>decay</code>	double radioactive half-life for solute ( )
<code>density_bulk</code>	double dry bulk density ( _b)
<code>porosity</code>	double porosity ( )
<code>tortuosity</code>	double tortuosity ( )
<code>n_terms</code>	integer number of terms for laplace inversion
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

**References**

Sudicky, E.A., Frind, E.O., Contaminant transport in fractured porous media: Analytical solutions for a system of parallel fractures, December 1982 <https://doi.org/10.1029/WR018i006p01634>

**See Also**

Other transport: [step\\_transport\\_fractures\\_heat\(\)](#), [step\\_transport\\_ogata\\_banks\(\)](#)

**Examples**

```
formula <- as.formula(~time+z+x)

dat <- as.data.frame(expand.grid(10^(3:8),
                               seq(0.0, 10, 1),
                               c(0.0)))

names(dat) <- c("time", "z", "x")

frec1 = recipe(formula = formula, data = dat) |>
  step_transport_fractures_solute(time = time,
                                distance_fracture = z,
                                distance_matrix = x) |>
  plate()
```

---

step\_transport\_ogata\_banks

*step\_transport\_ogata\_banks*

---

**Description**

Ogata, A., Banks, R.B., 1961. A solution of the differential equation of longitudinal dispersion in porous media. U. S. Geol. Surv. Prof. Pap. 411-A. 1-D, infinite source, uniform flow, constant parameters, decay, retardation

To have values match the excel sheet <https://www.civil.uwaterloo.ca/jrcraig/pdf/OgataBanks.xlsx> the decay coefficient needs to be scaled by the retardation coefficient.

Care must be taken so that input values do not lead to NaN. -Need to fix this.

1-D infinite source uniform flow constant parameters no decay no retardation

**Usage**

```
step_transport_ogata_banks(
  .rec,
  time,
  distance,
  concentration_initial = 1,
  velocity = 0.1,
  diffusion = 0.1,
  retardation = 1,
  decay = 0,
  role = "predictor",
  ...
)
```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	vector time
<code>distance</code>	vector x position
<code>concentration_initial</code>	double concentration
<code>velocity</code>	double velocity
<code>diffusion</code>	double diffusion coefficient
<code>retardation</code>	double retardation coefficient
<code>decay</code>	double decay coefficient
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

Ogata-Banks solution for time and distance pairs

**References**

Ogata, A., Banks, R.B., 1961. A solution of the differential equation of longitudinal dispersion in porous media. U. S. Geol. Surv. Prof. Pap. 411-A. 1-D, infinite source, uniform flow, constant parameters, decay, retardation

**See Also**

Other transport: [step\\_transport\\_fractures\\_heat\(\)](#), [step\\_transport\\_fractures\\_solute\(\)](#)

**Examples**

```

formula <- as.formula(y~x)
rows <- 100

dat <- data.frame(expand.grid(as.numeric(1:rows), as.numeric(1:10)))
names(dat) <- c('x', 'y')
frec1 = recipe(formula = formula, data = dat) |>
  step_transport_ogata_banks(time = x, distance = y) |>
  plate("dt")

```

---

```

step_vadose_weeks      step_vadose_weeks

```

---

**Description**

Weeks 1979 solution

**Usage**

```

step_vadose_weeks(
  .rec,
  time,
  air_diffusivity = 0.2,
  thickness = 40,
  precision = 1e-12,
  inverse = FALSE,
  role = "predictor",
  ...
)

```

**Arguments**

<code>.rec</code>	the R6 recipe object.
<code>time</code>	vector time elapsed time from start of pressure change
<code>air_diffusivity</code>	double vadose zone air diffusivity
<code>thickness</code>	double vadose zone thickness
<code>precision</code>	double stop the sum when precision is reached
<code>inverse</code>	double whether the response is invers
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

**Value**

an updated recipe

## References

Weeks, E.P., 1979. Barometric fluctuations in wells tapping deep unconfined aquifers. *Water Resources Research*, 15(5), pp.1167-1176.

## Examples

```
formula <- as.formula(y~x)

n <- 100
dat <- data.frame(x = as.numeric(1:n),
                 y = as.numeric(1:n))

frec1 = recipe(formula = formula, data = dat) |>
  step_vadose_weeks(time = x,
                   air_diffusivity = 0.8,
                   thickness = 5,
                   precision = 1e-12) |>
  plate()
```

---

step_varying	<i>step_varying</i>
--------------	---------------------

---

## Description

remove columns that only contain a single value.

## Usage

```
step_varying(.rec, terms, role = "predictor", ...)
```

## Arguments

<code>.rec</code>	the R6 recipe object.
<code>terms</code>	the unquoted names of the variables to use or a selector function. <code>terms</code> replaces the <code>'...'</code> of the <code>recipes</code> package but requires variables to be included within <code>'c()'</code> . For example to include variables <code>x</code> and <code>y</code> you would write <code>'c(x,y)'</code> in the <code>hydrorecipes</code> package.
<code>role</code>	character - the name of the role
<code>...</code>	additional arguments

## Value

an updated recipe

**Examples**

```

formula <- as.formula(y-x+z)
rows <- 1000
dat <- data.frame(x = rep(1, rows),
                  y = 1:rows,
                  z = rnorm(rows))

frec = recipe(formula = formula, data = dat) |>
  step_varying(c(x, y, z)) |>
  plate()

```

---

```
tidal_cooper_1965    tidal_cooper_1965
```

---

**Description**

Cooper Jr, H.H., Bredehoeft, J.D., Papadopoulos, I.S. and Bennett, R.R., 1965. The response of well-aquifer systems to seismic waves. *Journal of Geophysical Research*, 70(16), pp.3915-3926.

**Usage**

```

tidal_cooper_1965(
  frequency,
  storativity,
  transmissivity,
  thickness_aquifer,
  height_water,
  radius_well,
  radius_casing = radius_well,
  gravity = 9.80665
)

```

**Arguments**

<code>frequency</code>	the frequency of the response
<code>storativity</code>	layer storativity (L/L)
<code>transmissivity</code>	aquifer transmissivity (L*L/t)
<code>thickness_aquifer</code>	aquifer thickness
<code>height_water</code>	height of water in well
<code>radius_well</code>	well radius
<code>radius_casing</code>	casing radius
<code>gravity</code>	gravitational acceleration at well

**Value**

tidal response

**Examples**

```
data('hsieh_1987_fig_2_3')
storativity <- 1e-07
transmissivity <- 1e-03
radius_well <- 0.05
frequency <- 10^seq(-5, 2, by = 0.1)
tau <- 1 / frequency
cooper <- tidal_cooper_1965(frequency,
                           storativity,
                           transmissivity,
                           thickness_aquifer = 1,
                           height_water = 1,
                           radius_well)
plot(Mod(response)~dimensionless_frequency, cooper,
     type='l',
     log = 'x',
     xlim = c(1, 1000))
points(response~dimensionless_frequency,
       hsieh_1987_fig_2_3[variable=='gain' & S == storativity])

plot(unwrap(Arg(response)) * 180/pi~dimensionless_frequency, cooper,
     type='l',
     log = 'x',
     xlim = c(1, 1000),
     ylim = c(0, -90))
points(response~dimensionless_frequency,
       hsieh_1987_fig_2_3[variable=='phase' & S == storativity])
```

---

tidal\_hsieh\_1987

*tidal\_hsieh\_1987 Solution for estimating transmissivity and storativity from earth tides.*

---

**Description**

Hsieh, P.A., Bredehoeft, J.D. and Farr, J.M., 1987. Determination of aquifer transmissivity from Earth tide analysis. *Water resources research*, 23(10), pp.1824-1832.

**Usage**

```
tidal_hsieh_1987(
  frequency,
  storativity,
  transmissivity,
  radius_well,
```

```

    radius_casing = radius_well
  )

```

### Arguments

```

frequency      the frequency of the response
storativity    layer storativity (L/L)
transmissivity aquifer transmissivity (L*L/t)
radius_well    well radius
radius_casing  casing radius

```

### Value

tidal response

### Examples

```

data('hsieh_1987_fig_2_3')
storativity <- 1e-07
transmissivity <- 1e-03
radius_well <- 0.05
frequency <- 10^seq(-5, 2, by = 0.05)
tau <- 1 / frequency
hsieh <- tidal_hsieh_1987(frequency, storativity, transmissivity, radius_well)
plot(Mod(response)~dimensionless_frequency, hsieh,
     type='l',
     log = 'x',
     xlim = c(1, 1000))
points(response~dimensionless_frequency, hsieh_1987_fig_2_3[variable=='gain' & S == storativity])

plot(unwrap(Arg(response)) * 180/pi~dimensionless_frequency, hsieh,
     type='l',
     log = 'x',
     xlim = c(1, 1000),
     ylim = c(0, -90))
points(response~dimensionless_frequency, hsieh_1987_fig_2_3[variable=='phase' & S == storativity])

```

---

**unwrap**

*unwrap Removes the large phase shifts that can be associated with using Arg or atan2.*

---

### Description

unwrap Removes the large phase shifts that can be associated with using Arg or atan2.

**Usage**

```
unwrap(phase)
```

**Arguments**

`phase` the phase in radians

**Value**

unwrapped vector

---

<code>vadose_response</code>	<i>vadose_response</i>
------------------------------	------------------------

---

**Description**

weeks\_1979 1-D air diffusivity

**Usage**

```
vadose_response(time, air_diffusivity, thickness, precision, inverse)
```

**Arguments**

`time` numeric vector of elapsed times

`air_diffusivity` A numeric value of the unsaturated zone air diffusivity

`thickness` A numeric value of the unsaturated zone thickness

`precision` A numeric value of for the solution precision

`inverse` A logical value indicating if an inverse water level relationship is desired

**Value**

weeks 1979 model

**Examples**

```
vr <- vadose_response(time = 0:43200,
                      air_diffusivity = 0.20,
                      thickness = 40,
                      precision = 1e-10,
                      inverse = FALSE)
```

---

`window_blackman_harris`  
*window\_blackman\_harris*

---

### **Description**

Blackman-Harris window for FFT

### **Usage**

`window_blackman_harris(n)`

### **Arguments**

`n`                    length of the window vector (integer)

### **Value**

window

### **Examples**

`window_blackman_harris(100)`

---

`window_blackman_nuttall`  
*window\_blackman\_nuttall*

---

### **Description**

Blackman-Nuttall window for FFT

### **Usage**

`window_blackman_nuttall(n)`

### **Arguments**

`n`                    length of the window vector (integer)

### **Value**

window

**Examples**

```
window_blackman_nuttall(100)
```

---

```
window_first_deriv    window_first_deriv
```

---

**Description**

First derivative window for FFT

**Usage**

```
window_first_deriv(n, a0, a1, a2, a3)
```

**Arguments**

<code>n</code>	length of the window vector (integer)
<code>a0</code>	double coefficient
<code>a1</code>	double coefficient
<code>a2</code>	double coefficient
<code>a3</code>	double coefficient

**Value**

window

**Examples**

```
# nuttall window  
window_first_deriv(100, 0.355768, 0.487396, 0.144232, 0.012604)
```

---

window_hann	<i>window_hann</i>
-------------	--------------------

---

**Description**

Hann window for FFT.

**Usage**

window\_hann(n)

**Arguments**

n                    length of the window vector (integer)

**Value**

window of length n.

---

window_hann_cplx	<i>window_hann_cplx</i>
------------------	-------------------------

---

**Description**

Hann window for complex FFT.

**Usage**

window\_hann\_cplx(n)

**Arguments**

n                    length of the window vector (integer)

**Value**

window of length n.

`window_nuttall`      *window\_nuttall*

---

**Description**

Nuttall window for FFT

**Usage**

```
window_nuttall(n)
```

**Arguments**

`n`                    length of the window vector (integer)

**Value**

window

**Examples**

```
window_nuttall(100)
```

---

`window_rectangle`      *window\_rectangle*

---

**Description**

Rectangular window.

**Usage**

```
window_rectangle(n)
```

**Arguments**

`n`                    length of the window vector (integer)

**Value**

window of length `n`.

---

window_scale	<i>window_scale</i>
--------------	---------------------

---

**Description**

Scale factor for a window function.

**Usage**

```
window_scale(window, n_new, n_fft)
```

**Arguments**

window	the window function (numeric vector)
n_new	length of the padded series (integer)
n_fft	length of the input series (integer)

**Value**

window of length n.

---

window_tukey	<i>window_tukey</i>
--------------	---------------------

---

**Description**

Tukey window for FFT.

**Usage**

```
window_tukey(n, r)
```

**Arguments**

n	length of the window vector (integer)
r	percent on each side to taper (double)

**Value**

window of length n.

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