

# Package ‘asypDiag’

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**Type** Package

**Title** Diagnostic Tools for Asymptotic Theory

**Version** 0.3.2

**Description** Leveraging Monte Carlo simulations, this package provides tools for diagnosing regression models. It implements a parametric bootstrap framework to compute statistics, generates diagnostic envelopes to assess goodness-of-fit, and evaluates type I error control for Wald tests. By simulating data under the assumption that the model is true, it helps to identify model mis-specifications and enhances the reliability of the model inferences.

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---

concat_pvalues	<i>Concatenate AD_pvalues object</i>
----------------	--------------------------------------

---

### Description

Concatenate AD\_pvalues object

### Usage

```
concat_pvalues(ld_pvalues)
```

### Arguments

ld\_pvalues      list of elements of class AD\_pvalues.

### Value

Object of class AD\_pvalues

**Description**

Generates QQ-plot with simulated envelope residuals.

**Usage**

```
envelope(  
  model,  
  residual_fn = envelope_residual(model),  
  alpha = 0.05,  
  nsim = 100,  
  responses = NULL,  
  no_warnings = FALSE,  
  no_messages = FALSE,  
  converged_only = FALSE,  
  show_progress = TRUE,  
  plot.it = TRUE,  
  refit_fn = NULL,  
  ...  
)
```

**Arguments**

<code>model</code>	A model to generate responses and compute the observed residuals.
<code>residual_fn</code>	A function to calculate model residuals. The default is <code>envelope_residual()</code> for an absolute residual.
<code>alpha</code>	The significance level for constructing the envelope bounds. Defaults to 0.05.
<code>nsim</code>	The number of simulations to perform for envelope construction. Defaults to 100.
<code>responses</code>	An optional list of values to be used as response variables to refit the model.
<code>no_warnings</code>	If TRUE, ignore simulations that threw warnings.
<code>no_messages</code>	If TRUE, ignore simulations that shown messages.
<code>converged_only</code>	Use p-values from converged models only.
<code>show_progress</code>	Display a progress bar for the simulation iteration.
<code>plot.it</code>	Logical. Generate envelope plot.
<code>refit_fn</code>	Function to refit the model with new responses. If NULL, defaults to <code>get_refit(model, y, ...)</code> .
<code>...</code>	Extra arguments to <code>get_refit()</code>

## Details

Simulates new responses using `stats::simulate()` and refits the model for each vector of new responses using `get_refit()`. The function then computes residuals for each simulation, sorts them, and constructs envelope bands and a median line based on the quantiles of these residuals.

`refit_fn` is a function that supposedly compute the refit of model. Use this method if the default `get_refit()` doesn't work. If `refit_fn` is NULL, it's value is defined as `function(y, ...) get_refit(model, y, ...)`.

## Value

An object of class `AD_envelope`, which contains the following components:

**observed** A vector of observed quantiles from the model residuals.

**outside** A logical vector indicating whether each observation falls outside the constructed envelope bounds.

**lower** The lower bounds of the envelope for each observation.

**med** The median bounds of the envelope for each observation.

**upper** The upper bounds of the envelope for each observation.

## See Also

[get\\_refit](#), [simulate](#), [rstudent](#), [plot.AD\\_envelope](#), [parametric\\_bootstrap\(\)](#)

## Examples

```
fit <- lm(mpg ~ cyl, data = mtcars)

envelope(fit)

# Use pearson residuals, and plot it against the expected normal quantiles.
env_measures <- envelope(fit,
  residual_fn = function(x) residuals.lm(x, type = "pearson"), plot.it = FALSE
)
plot(env_measures, distribution = stats::qnorm, colors = c("gray", "black"))

## Using custom refit_fn
if (require("survival")) {
  fit <- survreg(Surv(futime, fustat) ~ ecog.ps + rx, ovarian,
    dist = "exponential"
  )
  fitted_rate <- 1 / fitted(fit)
  new_responses <- replicate(100, rexp(length(fitted_rate), fitted_rate), simplify = FALSE)
  refit_surv_ovarian <- function(.y) {
    survreg(Surv(.y, fustat) ~ ecog.ps + rx, ovarian, dist = "exponential")
  }
  env_measures <- envelope(fit,
    responses = new_responses,
    residual_fn = function(x) abs(residuals(x, type = "deviance")),
    refit_fn = refit_surv_ovarian, plot.it = FALSE
  )
}
```

```

)
# Absolute residuals are best shown with halfnormal quantiles
plot(env_measures, distribution = function(p) qnorm((1 + p) / 2))
}

```

---

envelope\_residual      *Recommended Residuals for Envelope Plots*

---

### Description

This function returns a function that computes residuals for envelope plots. These residuals are typically absolute values to be compared against the half-normal distribution.

### Usage

```

envelope_residual(object, ...)

## Default S3 method:
envelope_residual(object, ...)

## S3 method for class 'glm'
envelope_residual(object, ...)

## S3 method for class 'lm'
envelope_residual(object, ...)

```

### Arguments

`object`            An object for which model residuals can be extracted.  
`...`                Additional arguments passed to the residual function.

### Details

For objects of class `glm`, the default residuals are:

- Deviance residuals, except for poisson and binomial families.
- For poisson and binomial families it uses deletion residual adapted from [rstudent\(\)](#).

For objects of class `lm`, the default residuals is [rstudent\(\)](#).

For other classes, the default is [stats::residuals\(\)](#), meaning no specialized recommendation is currently provided.

### Value

A function that computes residuals from an object

---

get_converged	<i>Did the model converged</i>
---------------	--------------------------------

---

**Description**

Retrieves if the model converged.

**Usage**

```
get_converged(object)
```

```
## Default S3 method:  
get_converged(object)
```

**Arguments**

object            A model to check for convergence.

**Details**

The default method retrieves `object$converged`. For models of class `merMod` it verifies if the infinity norm of the Newton–Raphson step is less than 0.0001.

**Value**

A boolean value.

---

get_fixef	<i>Get fixed effects</i>
-----------	--------------------------

---

**Description**

Extracts the fixed effects coefficients from a model.

**Usage**

```
get_fixef(object)
```

```
## Default S3 method:  
get_fixef(object)
```

```
## S3 method for class 'merMod'  
get_fixef(object)
```

```
## S3 method for class 'glmmTMB'  
get_fixef(object)
```

**Arguments**

object            A model object for which fixed effects are to be retrieved.

**Details**

By default it calls `stats::coef()`. If the model class is `merMod` calls `fixef`.

**Value**

A numeric vector of fixed effects coefficients.

---

`get_model_response`     *Extract the Response Variable from a Model Object*

---

**Description**

This function is designed to extract the response variable from a fitted model object.

**Usage**

```
get_model_response(object)
```

**Arguments**

object            A fitted model from which the response variable should be extracted.

**Details**

The default method attempts to create a model frame using `model.frame()`. Any error encountered during this process is caught and results in a NULL return value.

**Value**

The response variable extracted from the model. If it couldn't be extracted, the function returns NULL.

**Examples**

```
lm_model <- lm(mpg ~ wt + cyl, data = mtcars)
response <- get_model_response(lm_model)
print(response)
```

---

get_refit	<i>Refit the Model</i>
-----------	------------------------

---

**Description**

Refit an existing regression model using a new response variable.

**Usage**

```
get_refit(object, newresp, ...)
```

**Arguments**

object	A model.
newresp	the new response, may be a vector or a matrix.
...	other arguments passed to refit or update.

**Details**

This generic method refit the object with the newresp. It should maintain the object properties by default, but some aspects of the default fitting method may be overwritten with the ...

**Value**

An updated model object, refitted using the new response variable.

**See Also**

[stats::update\(\)](#), [lme4::refit\(\)](#)

---

get_vcov	<i>Get covariance matrix</i>
----------	------------------------------

---

**Description**

Retrieves the covariance matrix for the fixed effects.

**Usage**

```
get_vcov(object)

## Default S3 method:
get_vcov(object)

## S3 method for class 'glmmTMB'
get_vcov(object)
```

**Arguments**

`object`            A model object for which the covariance matrix is to be retrieved.

**Details**

By default it calls `stats::vcov()` to retrieve the covariances.

**Value**

A covariance matrix of the model object.

---

parametric\_bootstrap    *Perform Parametric Bootstrap Simulations*

---

**Description**

This function performs parametric bootstrap simulations by generating new response values based on the fitted model, refitting the model on these new responses, and computing a user-defined statistic for each simulated model.

**Usage**

```
parametric_bootstrap(
  model,
  statistic,
  nsim,
  responses = NULL,
  refit_fn = NULL,
  show_progress = TRUE,
  simplify = TRUE,
  stat_hc = NULL,
  show_message_count = TRUE,
  show_warning_count = TRUE,
  show_not_converged_count = TRUE,
  ...
)
```

**Arguments**

`model`            A fitted model object that will be used to simulate responses.

`statistic`        A function that computes the desired statistic from the refitted model. It must take the refitted model as an argument.

`nsim`            The number of simulations to perform.

`responses`        An optional list of values to be used as response variables to refit the model.

`refit_fn`        Function to refit the model with new responses. If NULL, defaults to `get_refit(model, y, ...)`.

<code>show_progress</code>	Display a progress bar for the simulation iteration.
<code>simplify</code>	logical or character string; should the result be simplified to a vector, matrix or higher dimensional array if possible? If occurs any errors during the refit procedure, the results will be a list, regardless of the value of this argument.
<code>stat_hc</code>	A function that verifies if the computed statistic is correct. It should return nothing, just throw errors to halt execution.
<code>show_message_count</code>	Show total of captured messages from <code>refit_fn</code> as a message. It only shows if the number of messages is greater than 0.
<code>show_warning_count</code>	Show total of captured warnings from <code>refit_fn</code> as a warning. It only shows if the number of warnings is greater than 0.
<code>show_not_converged_count</code>	Show total of models that didn't converge as a warning. It only shows if the number of models that didn't converged is greater than 0.
<code>...</code>	Additional arguments to be passed to <code>refit_fn</code> .

### Details

This function implements a parametric bootstrap procedure. It generates new response values from the fitted model, refits the model for each simulated response, and computes a user-defined statistic on the refitted model. The refit function can be customized through the `refit_fn` argument.

If `show_progress` is TRUE, the progress of the simulations will be displayed using a progress bar from the `cli` package.

### Value

A list containing the following elements:

`result` A list of length `nsim` if `simplify` is FALSE. Otherwise an atomic vector or matrix or list of length `nsim`.

`responses` The list of simulated response values.

`simulation_warning` A logical vector indicating whether a warning occurred during model refitting for each simulation.

`converged` A logical vector indicating whether the model refit converged for each simulation.

### Examples

```
model <- lm(mpg ~ wt + hp, data = mtcars)
statistic <- function(model) coef(model)[["wt"]]
bootstrap_results <- parametric_bootstrap(model, statistic, nsim = 100)
wt_coefs <- Reduce(c, bootstrap_results$result)
summary(wt_coefs)
```

---

plot.AD_envelope	<i>Envelope Plot</i>
------------------	----------------------

---

## Description

Plot AD\_envelope

## Usage

```
## S3 method for class 'AD_envelope'
plot(
  x,
  colors = getOption("asymptDiag.plot.AD_envelope.colors"),
  xlab = "Expected quantiles",
  ylab = "Observed quantiles",
  distribution = function(p) stats::qnorm((1 + p)/2),
  ylim = base::range(c(x$observed, x$lower, x$upper), na.rm = TRUE, finite = TRUE),
  ...
)
```

## Arguments

x	AD_envelope object, usually the result of <a href="#">envelope()</a>
colors	Vector of length 2, with color for points outside and inside the envelope band, respectively.
xlab	The label for the x-axis.
ylab	The label for the y-axis.
distribution	quantile function for reference theoretical distribution.
ylim	the y limits of the plot.
...	extra arguments passed to <a href="#">graphics::plot</a>

## Details

Create envelope plot. The expected quantile, by default, is from a half-normal distribution, as the default residuals from [envelope\\_residual\(\)](#) are absolute. You may replace it with the distribution argument.

## Value

No return value, called for side effects

---

plot.AD\_pvalues      *Plot Empirical Cumulative Distribution Function (ECDF) of p-values*

---

### Description

This function creates several plots with the empirical cumulative distribution of the p-values obtained through simulation.

### Usage

```
## S3 method for class 'AD_pvalues'
plot(
  x,
  which = seq_len(length(x$test_coefficients) + 1),
  caption = as.list(paste("ECDF of", c(names(x$test_coefficients), "all coefficients"))),
  ks_test = TRUE,
  signif = c(0.01, 0.05, 0.1),
  discrepancy_tol = 0.1,
  plot_uniform = TRUE,
  uniform_legend = TRUE,
  converged_only = FALSE,
  no_warnings = FALSE,
  no_messages = FALSE,
  ylab = "Empirical cumulative distribution",
  xlab = "p-value",
  ...,
  ask = prod(graphics::par("mfcol")) < length(which) && grDevices::dev.interactive()
)
```

### Arguments

x	AD_pvalues object, usually the result of <code>simulate_wald_pvalues()</code>
which	A vector specifying the indices of coefficients to plot. If index is bigger than the number of coefficients it plots the joint p_value.
caption	A character vector or a list with caption for each plot. If it's a list, the list index must match the coefficient index used by which. If it's a vector, it's values are used in order.
ks_test	If TRUE inserts Kolmogorov-Smirnov p-value in the graphic.
signif	Points to verify discrepancy.
discrepancy_tol	Threshold to consider point discrepant.
plot_uniform	Logical. If TRUE, plot uniform distribution.
uniform_legend	Logical. If TRUE, a legend is added to the plot to distinguish between the p-value and U(0, 1) curves. Defaults to TRUE.
converged_only	Use p-values from converged models only.

no_warnings	If TRUE, ignore simulations that threw warnings.
no_messages	If TRUE, ignore simulations that shown messages.
ylab	The label for the y-axis. Defaults to "Empirical cumulative distribution".
xlab	The label for the x-axis. Defaults to "p-value".
...	extra arguments passed to <a href="#">graphics::plot</a>
ask	Logical. If TRUE, the user is prompted before each plot. Defaults to TRUE if in an interactive session and the number of plots is greater than the available space; otherwise, FALSE.

### Details

If the asymptotic approximation is valid the distribution of the p-values should be close to an uniform distribution. Discrepancies are highlighted, by default it verifies the significance on the most commonly used significance values are 0.01, 0.05 and 0.10.

The reported KS (Kolmogorov-Smirnov) test is the result of the "two-sided" `stats::ks.test()` function comparing the observed p-values distribution with the uniform. The test may reject the KS test due to few simulations, make sure that the lines shown in the plot are smooth before drawing any conclusions.

### Value

A vector of joint p-values for all coefficients.

### Examples

```
model <- lm(mpg ~ wt + hp, data = mtcars)
p_values_ld <- simulate_wald_pvalues(model, n_sim = 100)
plot(p_values_ld)
```

---

plot_cook	<i>Plot Cook's distances</i>
-----------	------------------------------

---

### Description

Plot Cook's distances

### Usage

```
plot_cook(
  model,
  n_highlights = 0,
  cut = FALSE,
  xlab = "Index",
  ylab = "Cook's distance",
  ...
)
```

**Arguments**

model	Model with <code>cooks.distance()</code> method
n_highlights	The number of observations with the highest Cook's distance to highlight on the plot. Defaults to 0 (no highlights).
cut	Logical. If TRUE, adds a cutoff line at the mean plus four times the standard deviation of Cook's distance. Defaults to FALSE.
xlab	The label for the x-axis. Defaults to "Index".
ylab	The label for the y-axis. Defaults to "Cook's distance".
...	Further arguments for <code>graphics::plot()</code>

**Value**

An invisible object representing Cook's distance values.

**See Also**

[cooks.distance](#), [plot](#)

**Examples**

```
fit <- lm(mpg ~ cyl, data = mtcars)

plot_cook(fit)
plot_cook(fit, n_highlights = 2)
```

---

plot\_ecdf\_pvalue      *Plot Empirical Cumulative Distribution Function (ECDF) of p-values*

---

**Description**

Plot Empirical Cumulative Distribution Function (ECDF) of p-values

**Usage**

```
plot_ecdf_pvalue(
  p_values,
  ks_test = TRUE,
  signif = c(0.01, 0.05, 0.1),
  discrepancy_tol = 0.1,
  plot_uniform = TRUE,
  uniform_legend = TRUE,
  main = "",
  ylab = "Empirical cumulative distribution",
  xlab = "p-value",
  ...
)
```

**Arguments**

p_values	vector of p-values
ks_test	If TRUE inserts Kolmogorov-Smirnov p-value in the graphic.
signif	Points to verify discrepancy.
discrepancy_tol	Threshold to consider point discrepant.
plot_uniform	Logical. If TRUE, plot uniform distribution.
uniform_legend	Logical. If TRUE, a legend is added to the plot to distinguish between the p-value and U(0, 1) curves. Defaults to TRUE.
main	main caption passed to <a href="#">plot</a>
ylab	The label for the y-axis. Defaults to "Empirical cumulative distribution".
xlab	The label for the x-axis. Defaults to "p-value".
...	extra arguments passed to <a href="#">graphics::plot</a>

**Value**

No return value, called for side effects

---

plot\_res\_vs\_linear\_predictor

*Plot Residuals against Linear Predictor*

---

**Description**

Plot Residuals against Linear Predictor

**Usage**

```
plot_res_vs_linear_predictor(
  model,
  residual_fn = stats::rstandard,
  xlab = "Linear Predictor",
  ylab = "Standardized deviance residuals",
  ...
)
```

**Arguments**

model	Model with methods <a href="#">predict()</a> or <a href="#">fitted()</a>
residual_fn	A function to calculate model residuals. The default is <code>stats::rstandard</code> .
xlab	The label for the x-axis. Defaults to "Linear Predictor".
ylab	The label for the y-axis. Defaults to "Standardized deviance residuals".
...	Extra arguments to <code>residual_fn</code> and <a href="#">plot()</a> .

**Details**

If the model was fitted using the `glm()` function, it will use the `predict()` method with `type = link`, otherwise, it will use the `fitted()` method.

**Value**

An invisible list containing the linear predictor (x) and standardized deviance residuals (y).

**Examples**

```
fit <- lm(mpg ~ cyl, data = mtcars)

plot_res_vs_linear_predictor(fit)
plot_res_vs_linear_predictor(fit, residual_fn = rstudent)
plot_res_vs_linear_predictor(fit, residual_fn = residuals)

glm_fit <- glm(cyl ~ mpg, family = poisson(), data = mtcars)

plot_res_vs_linear_predictor(glm_fit)
plot_res_vs_linear_predictor(glm_fit, type = "pearson")
```

---

refit\_model

*Refit Model*


---

**Description**

Refit a model with a new response.

**Usage**

```
refit_model(object, newresp, ...)
```

**Arguments**

<code>object</code>	A model.
<code>newresp</code>	the new response, may be a vector or a matrix.
<code>...</code>	other arguments passed to <code>refit</code> or <code>update</code> .

**Details**

This function uses `newresp` to refit `object` replacing its old response variable. If the class is `merMod` it uses `refit`, otherwise uses `stats::update()`.

The default method tries to update the model response using its `stats::model.frame()`, if it errors it tries to update the model by inserting the `newresp` directly into the object formula.

**Value**

A model with same class as object.

---

select_covariates	<i>Select covariates</i>
-------------------	--------------------------

---

**Description**

Select covariates

**Usage**

```
select_covariates(
  model,
  threshold = 0.15,
  direction = c("both", "backward", "forward"),
  addable_coefs = names(get_fixef(model)),
  measure_fn = function(x) summary(x)[["coefficients"]][, 4],
  measure_one_at_time = FALSE,
  minimize_only = FALSE,
  max_steps = 1000,
  return_step_results = FALSE,
  do_not_remove = c("(Intercept)"),
  ...
)
```

**Arguments**

model	A model with <code>stats::update()</code> , <code>stats::coef()</code> methods.
threshold	Value threshold to remove variable. It can be a fixed value or a function. The variable is removed if <code>measure_fn(model) &gt; threshold</code> and added if <code>measure_fn(model) &lt;= threshold</code> .
direction	The direction of variable selection. Options include "backward", "forward", or "both". Defaults to "both".
addable_coefs	A vector of coefficients that can be added during forward selection. Defaults to all coefficients in the model.
measure_fn	Function with model as argument and returns values to be used by threshold. It can also compare two models, where during forward step it calls <code>measure_fn(candidate_model, current_selected_model)</code> and during backward step it calls <code>measure_fn(current_selected_model, candidate_model)</code> . Defaults to the p-value from the summary of the coefficients.
measure_one_at_time	Boolean indicating to apply <code>measure_fn</code> to each variable individually during forward and backward steps. Set this option to TRUE if <code>measure_fn</code> returns an atomic value, for example if <code>measure_fn</code> is AIC.

`minimize_only` Logical indicating that during backward model update it should minimize the `measure_fn` instead of maximize it.

`max_steps` The maximum number of steps for the variable selection process. Defaults to 1000.

`return_step_results` Logical. If TRUE, the function returns a list containing the final fitted model and a log of the selection steps. Defaults to FALSE.

`do_not_remove` A character vector specifying variables that should not be removed during backward selection. Defaults to "(Intercept)".

... Extra arguments to `stats::update()`.

### Value

A fitted model with selected covariates based on the variable selection process. If `return_step_results` is TRUE, a list containing the final fitted model and a log of the selection steps is returned.

### Examples

```
model <- lm(mpg ~ ., data = mtcars)
select_covariates(model)

## measure_fn with two parameters

lrt <- function(model1, model2) {
  lrt_stat <- 2 * (logLik(model1)[1L] - logLik(model2)[1L])
  return(1 - pchisq(lrt_stat, 1))
}

select_covariates(model, measure_fn = lrt)

## AICc selection

AICc <- function(model) {
  loglike <- logLik(model)
  df <- attr(loglike, "df")
  nobs <- attr(loglike, "nobs")
  aic <- -2 * as.numeric(loglike) + 2 * df

  aicc <- aic + (2 * (df^2) + 2 * df) / (nobs - df - 1)

  return(aicc)
}

selection <- select_covariates(model,
  measure_fn = AICc,
  threshold = AICc,
  measure_one_at_time = TRUE,
  minimize_only = TRUE,
  direction = "both",
  data = mtcars
)
```

---

simulate\_wald\_pvalues *Generate Wald test P-Values with Monte Carlo Simulations*

---

### Description

This function performs Monte Carlo simulations to generate p-values for model coefficients by refitting the model with new simulated responses and computing the Wald test statistics for each simulation. It's standard behavior verify if the type I error from Wald tests are under control, considering the provided model as "true", i.e., the model assumptions are valid. It supports univariate and joint Wald tests, using chi-squared distributions to calculate p-values.

### Usage

```
simulate_wald_pvalues(
  model,
  nsim = 1000,
  responses = NULL,
  test_coefficients = NULL,
  refit_fn = NULL,
  coef_fn = get_fixef,
  vcov_fn = get_vcov,
  show_progress = TRUE,
  plot.it = TRUE,
  ...
)
```

### Arguments

model	A fitted model object that will be used to simulate responses.
nsim	The number of simulations to perform.
responses	An optional list of values to be used as response variables to refit the model.
test_coefficients	Numeric vector. A vector with values to be used to compute the test statistic. It should be the coefficients that was used to compute the fitted values of the response. If NULL defaults to <code>coef_fn(model)</code> .
refit_fn	Function to refit the model with new responses. If NULL, defaults to <code>get_refit(model, y, ...)</code> .
coef_fn	Function that retrieves the coefficients of the model.
vcov_fn	Function that computes the variance-covariance matrix for the models adjusted in the simulations.
show_progress	Display a progress bar for the simulation iteration.
plot.it	Logical. Generate ecdf plot for joint Wald test.
...	Additional arguments to be passed to <code>refit_fn</code> .

## Details

If `responses` is provided, the function refits the model with these new response vectors. Otherwise, it generates new responses with `stats::simulate()`.

For each new response calls `get_refit()` to generate a new model with the new response. It gets the fixed effects and the variance and covariance matrix with `get_fixef()` and `get_vcov()`.

Each simulated model is refitted using the specified `refit_fn` (or the default refit function) and the fixed effects coefficients and variance-covariance matrix are extracted using `coef_fn` and `vcov_fn`, respectively. The univariate Wald test is computed from the Wald statistic for each coefficient, while the joint Wald test uses the inverse variance-covariance matrix to compute a Wald statistic for the `test_coefficients`. P-values are calculated from a chi-squared distribution with appropriate degrees of freedom.

## Value

An object of class `AD_pvalues`, which contains the following components:

**test\_coefficients** Vector of coefficients being tested.

**pvalues\_matrix** Matrix of p-values where each column corresponds to a simulation and each row corresponds to a coefficient.

**pvalues\_joint** Vector containing the joint p-values obtained from each simulation.

**simulation\_fixef** List of fixed effect coefficient estimates from each simulation.

**simulation\_vcov** List of covariance matrices estimated from each simulation.

**simulation\_warning** Vector of boolean indicating if a simulation threw a warning.

**converged** Logical vector indicating whether model refitting converged for each simulation.

**responses** Simulated responses used for refitting the model.

## See Also

`plot.AD_pvalues()` for plotting.

## Examples

```
# from help("glm")
counts <- c(18, 17, 15, 20, 10, 20, 25, 13, 12)
outcome <- gl(3, 1, 9)
treatment <- gl(3, 3)
model <- glm(counts ~ outcome + treatment, family = poisson())
new_responses <- replicate(100, MASS::rnegbin(fitted.values(model), theta = 4.5), simplify = FALSE)
simulate_wald_pvalues(model, responses = new_responses, nsim = 100)

## Using custom refit_fn
if (require("survival")) {
  fit <- survreg(Surv(futime, fustat) ~ ecog.ps + rx, ovarian,
    dist = "exponential"
  )
  fitted_rate <- 1 / fitted(fit)
  new_responses <- replicate(100, rexp(length(fitted_rate), fitted_rate), simplify = FALSE)
```

```
refit_surv_ovarian <- function(.y) {  
  survreg(Surv(.y, fustat) ~ ecog.ps + rx, ovarian, dist = "exponential")  
}  
simulate_wald_pvalues(fit, responses = new_responses, refit_fn = refit_surv_ovarian)  
}
```

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