

Package ‘rstatix’

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Type Package

Title Pipe-Friendly Framework for Basic Statistical Tests

Version 1.1.0

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Description Provides a simple and intuitive pipe-friendly framework, coherent with the 'tidyverse' design philosophy, for performing basic statistical tests, including t-test, Wilcoxon test, ANOVA, Kruskal-Wallis and correlation analyses.
The output of each test is automatically transformed into a tidy data frame to facilitate visualization.
Additional functions are available for reshaping, reordering, manipulating and visualizing correlation matrix.
Functions are also included to facilitate the analysis of factorial experiments, including purely 'within-Ss' designs (repeated measures), purely 'between-Ss' designs, and mixed 'within-and-between-Ss' designs.
It's also possible to compute several effect size metrics, including ``eta squared" for ANOVA, ``Cohen's d" for t-test and 'Cramer V' for the association between categorical variables.
The package contains helper functions for identifying univariate and multivariate outliers, assessing normality and homogeneity of variances.

License GPL-2

Encoding UTF-8

Depends R (>= 3.3.0)

Imports stats, utils, tidyr (>= 1.0.0), purrr, broom (>= 0.7.4), rlang (>= 0.3.1), tibble (>= 2.1.3), dplyr (>= 0.7.1), magrittr, corplot, tidyselect (>= 1.2.0), car, generics (>= 0.0.2)

Suggests knitr, rmarkdown, ggpubr, graphics, emmeans, coin, boot, testthat (>= 3.1.7), spelling

VignetteBuilder knitr

URL <https://rpkgs.datanovia.com/rstatix/>

BugReports <https://github.com/kassambara/rstatix/issues>

Collate 'utilities.R' 'add_cld.R' 'add_significance.R'
 'adjust_pvalue.R' 'factorial_design.R'
 'utilities_two_sample_test.R' 'anova_summary.R' 'anova_test.R'
 'as_cor_mat.R' 'binom_test.R' 'box_m.R' 'posthoc_test.R'
 'check_test_assumptions.R' 'chisq_test.R' 'cliff_delta.R'
 'cochran_qtest.R' 'cohens_d.R' 't_test.R' 'dunn_test.R'
 'conover_test.R' 'cor_as_symbols.R' 'replace_triangle.R'
 'pull_triangle.R' 'cor_mark_significant.R' 'cor_mat.R'
 'cor_plot.R' 'cor_reorder.R' 'cor_reshape.R' 'cor_select.R'
 'cor_test.R' 'counts_to_cases.R' 'cramer_v.R' 'df.R' 'doo.R'
 'emmeans_test.R' 'dunnett_test.R' 'eta_squared.R' 'factors.R'
 'fisher_test.R' 'fligner_test.R' 'freq_table.R'
 'friedman_test.R' 'friedman_conover_test.R'
 'friedman_effsize.R' 'friedman_nemenyi_test.R'
 'games_howell_test.R' 'get_comparisons.R' 'get_mode.R'
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 'ks_test.R' 'levene_test.R' 'mahalanobis_distance.R'
 'make_clean_names.R' 'mcnemar_test.R' 'multinom_test.R'
 'omega_squared.R' 'outliers.R' 'p_value.R' 'prop_test.R'
 'prop_trend_test.R' 'reexports.R' 'remove_ns.R'
 'rstatix-programming.R' 'rstatix-references.R' 'sample_n_by.R'
 'shapiro_test.R' 'sign_test.R' 'tidy_glance.R' 'tukey_hsd.R'
 'utils-manova.R' 'utils-pipe.R' 'welch_anova_test.R'
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add_cld	<i>Compact Letter Display of All-Pairwise Comparisons</i>
---------	---

Description

Adds the **compact letter display** (CLD) to a pairwise comparison result. Groups that do *not* share a letter are significantly different. This is a convenient way to annotate plots (e.g. one letter per box/bar) after an all-pairwise post-hoc test such as `tukey_hsd()`, `dunn_test()`, `games_howell_test()`, `conover_test()`, `wilcox_test()` or `t_test()`.

The letters are computed with the insert-and-absorb algorithm (Piepho, 2004) using base R only, so no additional package is required (the results match `multcompView::multcomLetters()`).

Usage

```
add_cld(test, p.col = NULL, threshold = 0.05, reversed = FALSE, ...)
```

Arguments

- test an all-pairwise comparison result returned by an `rstatix` function (e.g. `tukey_hsd()`, `dunn_test()`, a pairwise `t_test()`/`wilcox_test()`, ...). Must contain the group1 and group2 columns and a p-value column.
- p.col character. The p-value column to threshold. If `NULL` (default), "p.adj" is used when present, otherwise "p".

threshold	the significance threshold (default 0.05). Comparisons with a p-value below threshold are treated as significant; comparisons with a missing (NA) p-value are treated as non-significant.
reversed	logical. If TRUE, reverses the order in which the letters are assigned (so that, with groups ordered by increasing level, the later groups receive the earlier letters). Default is FALSE.
...	not used.

Value

a tibble with one row per group and the following columns: any grouping variables (for a grouped test), .y. (the outcome variable, when present), group (the group level) and cld (the compact letter display). Groups sharing a letter are not significantly different.

References

Piepho, H.-P. (2004) An Algorithm for a Letter-Based Representation of All-Pairwise Comparisons. *Journal of Computational and Graphical Statistics*, 13(2), 456-466.

See Also

[tukey_hsd](#), [dunn_test](#), [games_howell_test](#), [add_significance](#)

Examples

```
# Tukey HSD post-hoc, then compact letter display
res <- ToothGrowth %>%
  mutate(dose = factor(dose)) %>%
  tukey_hsd(len ~ dose)
res %>% add_cld()

# Works on rank-based post-hocs too
ToothGrowth %>% dunn_test(len ~ dose) %>% add_cld()

# Grouped pairwise test -> one CLD per group
ToothGrowth %>%
  mutate(dose = factor(dose)) %>%
  group_by(supp) %>%
  tukey_hsd(len ~ dose) %>%
  add_cld()
```

add_significance	<i>Add P-value Significance Symbols</i>
------------------	---

Description

Add p-value significance symbols into a data frame.

Usage

```
add_significance(
  data,
  p.col = NULL,
  output.col = NULL,
  cutpoints = c(0, 1e-04, 0.001, 0.01, 0.05, 1),
  symbols = c("****", "***", "**", "*", "ns")
)
```

Arguments

<code>data</code>	a data frame containing a p-value column.
<code>p.col</code>	column name containing p-values.
<code>output.col</code>	the output column name to hold the adjusted p-values.
<code>cutpoints</code>	numeric vector used for intervals.
<code>symbols</code>	character vector, one shorter than cutpoints, used as significance symbols.

Value

a data frame

Examples

```
# Perform pairwise comparisons and adjust p-values
ToothGrowth %>%
  t_test(len ~ dose) %>%
  adjust_pvalue() %>%
  add_significance("p.adj")
```

adjust_pvalue

Adjust P-values for Multiple Comparisons

Description

A pipe-friendly function to add an adjusted p-value column into a data frame. Supports grouped data.

Usage

```
adjust_pvalue(data, p.col = NULL, output.col = NULL, method = "holm")
```

Arguments

data	a data frame containing a p-value column
p.col	column name containing p-values
output.col	the output column name to hold the adjusted p-values
method	method for adjusting p values (see p.adjust). Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use p.adjust.method = "none".

Details

For **grouped data** (and, equivalently, when a test is run on data grouped with `dplyr::group_by()` using an in-test `p.adjust.method`), the p-value adjustment is computed **within each group separately**, not across all groups. If you instead want a single family of comparisons adjusted across all groups, run the test without adjustment (`p.adjust.method = "none"`) and then call `adjust_pvalue()` on the combined result (see the grouped example below).

Value

a data frame

Examples

```
# Perform pairwise comparisons and adjust p-values
ToothGrowth %>%
  t_test(len ~ dose) %>%
  adjust_pvalue()

# Grouped data: adjustment within vs across groups
# Per-group adjustment (within each supp level):
ToothGrowth %>%
  group_by(supp) %>%
  t_test(len ~ dose)                # in-test holm, adjusted within each group

# One family across ALL comparisons (all groups together):
ToothGrowth %>%
  group_by(supp) %>%
  t_test(len ~ dose, p.adjust.method = "none") %>%
  adjust_pvalue(method = "holm")
```

 anova_summary

Create Nice Summary Tables of ANOVA Results

Description

Create beautiful summary tables of ANOVA test results obtained from either [Anova\(\)](#) or [aov\(\)](#).

The results include ANOVA table, generalized effect size and some assumption checks.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
anova_summary(
  object,
  effect.size = "ges",
  detailed = FALSE,
  observed = NULL,
  ci = NULL
)
```

Arguments

<code>object</code>	an object of returned by either <code>Anova()</code> , or <code>aov()</code> .
<code>effect.size</code>	the effect size to compute and to show in the ANOVA results. Allowed values can be either "ges" (generalized eta squared) or "pes" (partial eta squared) or both. Default is "ges".
<code>detailed</code>	If TRUE, returns extra information (sums of squares columns, intercept row, etc.) in the ANOVA table.
<code>observed</code>	Variables that are observed (i.e, measured) as compared to experimentally manipulated. The default effect size reported (generalized eta-squared) requires correct specification of the observed variables.
<code>ci</code>	confidence level for a confidence interval on partial eta squared. If a number between 0 and 1 (e.g. 0.95) and <code>effect.size</code> includes "pes", adds <code>conf.low/conf.high</code> columns computed from the noncentral F distribution. Default NULL (no interval).

Value

return an object of class `anova_test` a data frame containing the ANOVA table for independent measures ANOVA. However, for repeated/mixed measures ANOVA, it is a list containing the following components are returned:

- **ANOVA:** a data frame containing ANOVA results
- **Mauchly's Test for Sphericity:** If any within-Ss variables with more than 2 levels are present, a data frame containing the results of Mauchly's test for Sphericity. Only reported for effects that have more than 2 levels because sphericity necessarily holds for effects with only 2 levels.
- **Sphericity Corrections:** If any within-Ss variables are present, a data frame containing the Greenhouse-Geisser and Huynh-Feldt epsilon values, and corresponding corrected p-values.

The **returned object might have an attribute** called `args` if you compute ANOVA using the function `anova_test()`. The attribute `args` is a list holding the arguments used to fit the ANOVA model, including: `data`, `dv`, `within`, `between`, `type`, `model`, etc.

The following abbreviations are used in the different results tables:

- **DFn** Degrees of Freedom in the numerator (i.e. DF effect).
- **DFd** Degrees of Freedom in the denominator (i.e., DF error).
- **SSn** Sum of Squares in the numerator (i.e., SS effect).

- SSd Sum of Squares in the denominator (i.e.,SS error).
- F F-value.
- p p-value (probability of the data given the null hypothesis).
- p<.05 Highlights p-values less than the traditional alpha level of .05.
- ges Generalized Eta-Squared measure of effect size.
- GGe Greenhouse-Geisser epsilon.
- p[GGe] p-value after correction using Greenhouse-Geisser epsilon.
- p[GGe]<.05 Highlights p-values (after correction using Greenhouse-Geisser epsilon) less than the traditional alpha level of .05.
- HFe Huynh-Feldt epsilon.
- p[HFe] p-value after correction using Huynh-Feldt epsilon.
- p[HFe]<.05 Highlights p-values (after correction using Huynh-Feldt epsilon) less than the traditional alpha level of .05.
- W Mauchly's W statistic

Author(s)

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See Also

[anova_test\(\)](#), [factorial_design\(\)](#) The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth
df$dose <- as.factor(df$dose)

# Independent measures ANOVA
#::::::::::::::::::::::::::::::::::::
# Compute ANOVA and display the summary
res.anova <- Anova(lm(len ~ dose*supp, data = df))
anova_summary(res.anova)

# Display both SSn and SSd using detailed = TRUE
# Show generalized eta squared using effect.size = "ges"
anova_summary(res.anova, detailed = TRUE, effect.size = "ges")

# Show partial eta squared using effect.size = "pes"
anova_summary(res.anova, detailed = TRUE, effect.size = "pes")

# Repeated measures designs using car::Anova()
#::::::::::::::::::::::::::::::::::::
# Prepare the data
df$id <- as.factor(rep(1:10, 6)) # Add individuals ids
```

```

head(df)

# Easily perform repeated measures ANOVA using the car package
design <- factorial_design(df, dv = len, wid = id, within = c(supp, dose))
res.anova <- Anova(design$model, idata = design$idata, idesign = design$idesign, type = 3)
anova_summary(res.anova)

# Repeated measures designs using stats::Aov()
#::::::::::::::::::::::::::::::::::::::::::::::::::
res.anova <- aov(len ~ dose*supp + Error(id/(supp*dose)), data = df)
anova_summary(res.anova)

```

anova_test

Anova Test

Description

Provides a pipe-friendly framework to perform different types of ANOVA tests, including:

- **Independent measures ANOVA:** between-Subjects designs,
- **Repeated measures ANOVA:** within-Subjects designs
- **Mixed ANOVA:** Mixed within within- and between-Subjects designs, also known as split-plot ANOVA and
- **ANCOVA: Analysis of Covariance.**

The function is an easy to use wrapper around [Anova\(\)](#) and [aov\(\)](#). It makes ANOVA computation handy in R and It's highly flexible: can support model and formula as input. Variables can be also specified as character vector using the arguments `dv`, `wid`, `between`, `within`, `covariate`.

The results include ANOVA table, generalized effect size and some assumption checks.

Usage

```

anova_test(
  data,
  formula,
  dv,
  wid,
  between,
  within,
  covariate,
  type = NULL,
  effect.size = "ges",
  error = NULL,
  white.adjust = FALSE,
  observed = NULL,
  detailed = FALSE,
  ci = NULL

```

```

)

get_anova_table(x, correction = c("auto", "GG", "HF", "none"))

## S3 method for class 'anova_test'
print(x, ...)

## S3 method for class 'anova_test'
plot(x, ...)

```

Arguments

data	a data.frame or a model to be analyzed.
formula	a formula specifying the ANOVA model similar to aov. Can be of the form $y \sim \text{group}$ where y is a numeric variable giving the data values and group is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code>. Examples of supported formula include: • Between-Ss ANOVA (independent measures ANOVA): $y \sim b1*b2$ • Within-Ss ANOVA (repeated measures ANOVA): $y \sim w1*w2 + \text{Error}(\text{id}/(w1*w2))$ • Mixed ANOVA: $y \sim b1*b2*w1 + \text{Error}(\text{id}/w1)$ If the formula doesn't contain any within vars, a linear model is directly fitted and passed to the ANOVA function. For repeated designs, the ANOVA variables are parsed from the formula.
dv	(numeric) dependent variable name.
wid	(factor) column name containing individuals/subjects identifier. Should be unique per individual.
between	(optional) between-subject factor variables.
within	(optional) within-subjects factor variables
covariate	(optional) covariate names (for ANCOVA)
type	the type of sums of squares for ANOVA. Allowed values are either 1, 2 or 3. <code>type = 2</code> is the recommended default because it yields identical ANOVA results to <code>type = 1</code> when data are balanced, while also producing various assumption tests where appropriate. When the data are unbalanced, <code>type = 3</code> is the choice used by popular commercial softwares including SPSS. Default when type is not specified: the chosen default depends on the interface (see Note), so for unbalanced designs the two interfaces can differ. To get reproducible, interface-independent results, set type explicitly (e.g. <code>type = 3</code> for an unbalanced factorial design with interactions).
effect.size	the effect size to compute and to show in the ANOVA results. Allowed values can be either "ges" (generalized eta squared) or "pes" (partial eta squared) or both. Default is "ges".
error	(optional) for a linear model, an lm model object from which the overall error sum of squares and degrees of freedom are to be calculated. Read more in Anova() documentation.

<code>white.adjust</code>	Default is FALSE. If TRUE, heteroscedasticity correction is applied to the coefficient of covariance matrix. Used only for independent measures ANOVA.
<code>observed</code>	Variables that are observed (i.e, measured) as compared to experimentally manipulated. The default effect size reported (generalized eta-squared) requires correct specification of the observed variables.
<code>detailed</code>	If TRUE, returns extra information (sums of squares columns, intercept row, etc.) in the ANOVA table.
<code>ci</code>	confidence level for a confidence interval on the effect size. If a number between 0 and 1 (e.g. 0.95), two columns <code>conf.low</code> and <code>conf.high</code> are added giving the confidence interval for partial eta squared (<code>effect.size</code> must include "pes"). The interval is computed from the noncentral F distribution (Steiger, 2004) in base R. The interval is two-sided, and matches <code>effectsize::eta_squared(partial = TRUE, ci = , alternative = "two.sided")</code> ; that function defaults to a one-sided interval, whose upper bound is 1. No interval is provided for generalized eta squared ("ges"), which has no standard closed-form interval. Default is NULL (no interval; output unchanged).
<code>x</code>	an object of class <code>anova_test</code>
<code>correction</code>	character. Used only in repeated measures ANOVA test to specify which correction of the degrees of freedom should be reported for the within-subject factors. Possible values are: • "GG": applies Greenhouse-Geisser correction to all within-subjects factors even if the assumption of sphericity is met (i.e., Mauchly's test is not significant, $p > 0.05$). • "HF": applies Huynh-Feldt correction to all within-subjects factors even if the assumption of sphericity is met, • "none": returns the ANOVA table without any correction and • "auto": apply automatically GG correction to only within-subjects factors violating the sphericity assumption (i.e., Mauchly's test p-value is significant, $p \leq 0.05$).
<code>...</code>	additional arguments

Details

Contrasts. By default, R uses treatment contrasts (`contr.treatment`), where each factor level is compared to the first level used as baseline; the current setting can be checked with `options('contrasts')`.

How `anova_test()` handles contrasts depends on the interface you use:

- When you use the **formula interface** (or the `dv + between/within` arguments), `anova_test()` fits the model internally with `options(contrasts = c('contr.sum', 'contr.poly'))`, restoring your global option afterwards. This gives orthogonal contrasts, where every level is compared to the overall mean, and type-III results that match the most commonly used commercial softwares, like SPSS.
- When you instead pass a **pre-fitted model** (`lm()` or `aov()`), `anova_test()` does not change its contrasts: the model keeps whatever contrasts were in effect when it was fitted (R's default `contr.treatment` unless you set otherwise). Fitting with the default treatment contrasts and then requesting `type = 3` can therefore give different results from the formula interface.

To reproduce the formula-interface (SPSS) result from a pre-fitted model, or to obtain the same result with `car::Anova()` directly, set `options(contrasts = c('contr.sum', 'contr.poly'))` *before* fitting the model and use `type = 3`.

Value

return an object of class `anova_test` a data frame containing the ANOVA table for independent measures ANOVA.

However, for repeated/mixed measures ANOVA, a list containing the following components are returned: ANOVA table, Mauchly's Test for Sphericity, Sphericity Corrections. These table are described more in the documentation of the function `anova_summary()`.

The **returned object has an attribute** called `args`, which is a list holding the arguments used to fit the ANOVA model, including: `data`, `dv`, `within`, `between`, `type`, `model`, etc.

Functions

- `anova_test()`: perform anova test
- `get_anova_table()`: extract anova table from an object of class `anova_test`. When within-subject factors are present, either sphericity corrected or uncorrected degrees of freedom can be reported.

Note

Default sums-of-squares type differs between the two interfaces for unbalanced designs. When `type` is not supplied:

- the **formula** interface (`anova_test(data, y ~ a*b)`) uses **type II** for between-subjects designs, regardless of balance;
- the `dv=/between=/within=` interface uses type II for *balanced* between-subjects designs but switches to **type III** for *unbalanced* between-subjects designs with more than one factor (both interfaces use type III for repeated-measures designs).

For a **balanced** design the SS types coincide, so the two interfaces agree. For an **unbalanced** factorial design they can give different main-effect F/p values unless you **pass type explicitly** — which is recommended for reproducibility. Example:

```
d <- mtcars %>% dplyr::mutate(cyl = factor(cyl), am = factor(am))
# differ (unbalanced): formula -> type II, dv/between -> type III
d %>% anova_test(mpg ~ cyl * am)
d %>% anova_test(dv = mpg, between = c(cyl, am))
# agree once type is explicit:
d %>% anova_test(mpg ~ cyl * am, type = 3)
d %>% anova_test(dv = mpg, between = c(cyl, am), type = 3)
```

Author(s)

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See Also

[anova_summary\(\)](#), [factorial_design\(\)](#) The Datanovia tutorials: [One-Way ANOVA in R](#), [Repeated Measures ANOVA in R](#), [Mixed ANOVA in R](#), [ANCOVA in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# One-way ANOVA test
#::::::::::::::::::::::::::::::::::::
df %>% anova_test(len ~ dose)

# Grouped One-way ANOVA test
#::::::::::::::::::::::::::::::::::::
df %>%
  group_by(supp) %>%
  anova_test(len ~ dose)

# Two-way ANOVA test
#::::::::::::::::::::::::::::::::::::
df %>% anova_test(len ~ supp*dose)

# Two-way repeated measures ANOVA
#::::::::::::::::::::::::::::::::::::
df$id <- rep(1:10, 6) # Add individuals id
# Use formula

df %>% anova_test(len ~ supp*dose + Error(id/(supp*dose)))

# or use character vector
df %>% anova_test(dv = len, wid = id, within = c(supp, dose))

# Extract ANOVA table and apply correction
#::::::::::::::::::::::::::::::::::::
res.aov <- df %>% anova_test(dv = len, wid = id, within = c(supp, dose))
get_anova_table(res.aov, correction = "GG")

# Use model as arguments
#::::::::::::::::::::::::::::::::::::
.my.model <- lm(yield ~ block + N*P*K, npk)
anova_test(.my.model)
```

as_cor_mat	<i>Convert a Correlation Test Data Frame into a Correlation Matrix</i>
------------	--

Description

Convert a correlation test data frame, returned by the `cor_test()`, into a correlation matrix format.

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

```
as_cor_mat(x)
```

Arguments

x an object of class `cor_test`.

Value

Returns a data frame containing the matrix of the correlation coefficients. The output has an attribute named "pvalue", which contains the matrix of the correlation test p-values.

See Also

`cor_mat()`, `cor_test()` The Datanovia tutorial: **Correlation Matrix in R**.

Examples

[illegible]

binom_test

*Exact Binomial Test***Description**

Performs exact binomial test and pairwise comparisons following a significant exact multinomial test. Wrapper around the R base function `link[stats]{binom.test}()` that returns a data frame as a result.

Usage

```
binom_test(
  x,
  n,
  p = 0.5,
  alternative = "two.sided",
  conf.level = 0.95,
  detailed = FALSE
)

pairwise_binom_test(
  x,
  p.adjust.method = "holm",
  alternative = "two.sided",
  conf.level = 0.95
)

pairwise_binom_test_against_p(
  x,
  p = rep(1/length(x), length(x)),
  p.adjust.method = "holm",
  alternative = "two.sided",
  conf.level = 0.95
)
```

Arguments

<code>x</code>	numeric vector containing the counts.
<code>n</code>	number of trials; ignored if <code>x</code> has length 2.
<code>p</code>	a vector of probabilities of success. The length of <code>p</code> must be the same as the number of groups specified by <code>x</code> , and its elements must be greater than 0 and less than 1.
<code>alternative</code>	indicates the alternative hypothesis and must be one of "two.sided", "greater" or "less". You can specify just the initial letter.
<code>conf.level</code>	confidence level for the returned confidence interval.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.

`p.adjust.method`

method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use `p.adjust.method = "none"`.

Value

return a data frame containing the p-value and its significance. with some the following columns:

- `group`, `group1`, `group2`: the categories or groups being compared.
- `statistic`: the number of successes.
- `parameter`: the number of trials.
- `p`: p-value of the test.
- `p.adj`: the adjusted p-value.
- `method`: the used statistical test.
- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: the estimated probability of success.
- `alternative`: a character string describing the alternative hypothesis.
- `conf.low`, `conf.high`: Lower and upper bound on a confidence interval for the probability of success.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `binom_test()`: performs exact binomial test. Wrapper around the R base function `binom.test` that returns a dataframe as a result.
- `pairwise_binom_test()`: performs pairwise comparisons (binomial test) following a significant exact multinomial test.
- `pairwise_binom_test_against_p()`: performs pairwise comparisons (binomial test) following a significant exact multinomial test for given probabilities.

See Also

[multinom_test](#)

Examples

```
# Exact binomial test
#####
# Data: 160 mice with cancer including 95 male and 65 female
# Q1: Does cancer affect more males than females?
binom_test(x = 95, n = 160)
# => yes, there are a significant difference
```

```

# Q2: compare the observed proportion of males
# to an expected proportion (p = 3/5)
binom_test(x = 95, n = 160, p = 3/5)
# => there are no significant difference

# Multinomial test
#####
# Data
tulip <- c(red = 81, yellow = 50, white = 27)
# Question 1: are the color equally common ?
# this is a test of homogeneity
res <- multinom_test(tulip)
res
attr(res, "descriptives")

# Pairwise comparisons between groups
pairwise_binom_test(tulip, p.adjust.method = "bonferroni")

# Question 2: comparing observed to expected proportions
# this is a goodness-of-fit test
expected.p <- c(red = 0.5, yellow = 0.33, white = 0.17)
res <- multinom_test(tulip, expected.p)
res
attr(res, "descriptives")

# Pairwise comparisons against a given probabilities
pairwise_binom_test_against_p(tulip, expected.p)

```

box_m

Box's M-test for Homogeneity of Covariance Matrices

Description

Performs the Box's M-test for homogeneity of covariance matrices obtained from multivariate normal data according to one grouping variable. The test is based on the chi-square approximation.

See the Datanovia tutorial [Homogeneity of Variance Test in R](#) for a worked walkthrough.

Usage

```
box_m(data, group)
```

Arguments

data	a numeric data.frame or matrix containing n observations of p variables; it is expected that $n > p$.
group	a vector of length n containing the class of each observation; it is usually a factor.

Value

A data frame containing the following components:

statistic	an approximated value of the chi-square distribution.
parameter	the degrees of freedom related of the test statistic in this case that it follows a Chi-square distribution.
p.value	the p-value of the test.
method	the character string "Box's M-test for Homogeneity of Covariance Matrices".

See Also

The Datanovia tutorial: [Homogeneity of Variance Test in R](#).

Examples

```
data(iris)
box_m(iris[, -5], iris[, 5])
```

check_test_assumptions

Check One-Way Assumptions and Recommend the Test

Description

For a one-way, independent-groups design (outcome ~ group), check the two assumptions that decide which family of tests is appropriate — normality (Shapiro-Wilk, per group) and homogeneity of variance (Levene) — and return the verdicts together with the recommended omnibus and post-hoc test:

- each group normal **and** equal variances: `anova_test() + tukey_hsd()`;
- each group normal **but** unequal variances: `welch_anova_test() + games_howell_test()`;
- at least one group not normal: `kruskal_test() + dunn_test()`.

The result is a tidy one-row tibble, so the same single assumption check can drive both the omnibus and the post-hoc coherently — run the recommended omnibus, then pass the result to `posthoc_test()` via its `.assumptions` argument to avoid re-checking.

A note on choosing a test from the data. Selecting the test by first testing its assumptions on the same data is convenient but has a known cost: the assumption gate is least reliable exactly when it matters (Shapiro-Wilk has little power at small *n* and rejects trivial departures at large *n*), and conditioning the choice on it makes the p-value of the test finally run no longer the exact nominal quantity. A common alternative is to skip the gate and use a robust method unconditionally — Welch ANOVA with Games-Howell (which reduce to the classic result when variances are equal) or a rank-based test. Treat this recommendation as guidance, not a substitute for judgement.

See the Datanovia tutorial [Statistical Tests and Assumptions in R](#) for a worked walkthrough.

Usage

```
check_test_assumptions(data, formula, significance = 0.05)
```

Arguments

data	a data frame containing the variables in the formula.
formula	a formula of the form $x \sim \text{group}$ where x is a numeric outcome and group is a factor with two or more levels.
significance	the significance level used to judge the Shapiro-Wilk and Levene tests. Default is 0.05.

Value

a one-row tibble with the columns `.y.` (the outcome), `normality.p` (the smallest Shapiro-Wilk p across groups), `homogeneity.p` (Levene's p), the logical verdicts `normal` and `equal.variance`, the significance used, and the recommended omnibus and posthoc test names.

See Also

`posthoc_test()`, `shapiro_test()`, `levene_test()`. The Datanovia tutorial: [Statistical Tests and Assumptions in R](#).

Examples

```
df <- ToothGrowth
df$dose <- as.factor(df$dose)
df %>% check_test_assumptions(len ~ dose)
```

chisq_test

Chi-squared Test for Count Data

Description

Performs chi-squared tests, including goodness-of-fit, homogeneity and independence tests.

`chisq_test()` also accepts a pipe-friendly data-frame interface for the test of independence between two categorical variables: pass a data frame as `x` and the two columns either positionally (`data %>% chisq_test(var1, var2)`) or via `vars` (`data %>% chisq_test(vars = c("var1", "var2"))`). The contingency table is built internally. Note that in the positional form the second column occupies the correct argument slot, so use the `vars` form (or the table interface) if you need to set `correct/simulate.p.value`.

See the Datanovia tutorial [Chi-Square Test of Independence in R](#) for a worked walkthrough.

Usage

```

chisq_test(
  x,
  y = NULL,
  correct = TRUE,
  p = rep(1/length(x), length(x)),
  rescale.p = FALSE,
  simulate.p.value = FALSE,
  B = 2000,
  vars = NULL
)

pairwise_chisq_gof_test(x, p.adjust.method = "holm", ...)

pairwise_chisq_test_against_p(
  x,
  p = rep(1/length(x), length(x)),
  p.adjust.method = "holm",
  ...
)

chisq_descriptives(res.chisq)

expected_freq(res.chisq)

observed_freq(res.chisq)

pearson_residuals(res.chisq)

std_residuals(res.chisq)

```

Arguments

<code>x</code>	a numeric vector or matrix. <code>x</code> and <code>y</code> can also both be factors.
<code>y</code>	a numeric vector; ignored if <code>x</code> is a matrix. If <code>x</code> is a factor, <code>y</code> should be a factor of the same length.
<code>correct</code>	a logical indicating whether to apply continuity correction when computing the test statistic for 2 by 2 tables: one half is subtracted from all $ O - E $ differences; however, the correction will not be bigger than the differences themselves. No correction is done if <code>simulate.p.value = TRUE</code> .
<code>p</code>	a vector of probabilities of the same length as <code>x</code> . An error is given if any entry of <code>p</code> is negative.
<code>rescale.p</code>	a logical scalar; if <code>TRUE</code> then <code>p</code> is rescaled (if necessary) to sum to 1. If <code>rescale.p</code> is <code>FALSE</code> , and <code>p</code> does not sum to 1, an error is given.
<code>simulate.p.value</code>	a logical indicating whether to compute p-values by Monte Carlo simulation.
<code>B</code>	an integer specifying the number of replicates used in the Monte Carlo test.

<code>vars</code>	optional character vector of length two giving the names of two columns in the data frame <code>x</code> to cross-tabulate for a test of independence. An alternative to passing the two columns positionally.
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>...</code>	other arguments passed to the function <code>{chisq_test}()</code> .
<code>res.chisq</code>	an object of class <code>chisq_test</code> .

Value

return a data frame with some the following columns:

- `n`: the number of participants.
- `group`, `group1`, `group2`: the categories or groups being compared.
- `statistic`: the value of Pearson's chi-squared test statistic.
- `df`: the degrees of freedom of the approximate chi-squared distribution of the test statistic. NA if the p-value is computed by Monte Carlo simulation.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the used statistical test.
- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `observed`: observed counts.
- `expected`: the expected counts under the null hypothesis.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `chisq_test()`: performs chi-square tests including goodness-of-fit, homogeneity and independence tests.
- `pairwise_chisq_gof_test()`: perform pairwise comparisons between groups following a global chi-square goodness of fit test.
- `pairwise_chisq_test_against_p()`: perform pairwise comparisons after a global chi-squared test for given probabilities. For each group, the observed and the expected proportions are shown. Each group is compared to the sum of all others.
- `chisq_descriptives()`: returns the descriptive statistics of the chi-square test. These include, observed and expected frequencies, proportions, residuals and standardized residuals. Only available for a single (ungrouped) `chisq_test()` result.
- `expected_freq()`: returns the expected counts from the chi-square test result.
- `observed_freq()`: returns the observed counts from the chi-square test result.
- `pearson_residuals()`: returns the Pearson residuals, $(\text{observed} - \text{expected}) / \sqrt{\text{expected}}$.
- `std_residuals()`: returns the standardized residuals

See Also

The Datanovia tutorial: [Chi-Square Test of Independence in R](#), [Chi-Square Goodness-of-Fit Test in R](#).

Examples

```
# Chi-square goodness of fit test
#####
tulip <- c(red = 81, yellow = 50, white = 27)
# Q1: Are the colors equally common?
chisq_test(tulip)
pairwise_chisq_gof_test(tulip)
# Q2: comparing observed to expected proportions
chisq_test(tulip, p = c(1/2, 1/3, 1/6))
pairwise_chisq_test_against_p(tulip, p = c(0.5, 0.33, 0.17))

# Homogeneity of proportions between groups
#####
# Data: Titanic
xtab <- as.table(rbind(
  c(203, 118, 178, 212),
  c(122, 167, 528, 673)
))
dimnames(xtab) <- list(
  Survived = c("Yes", "No"),
  Class = c("1st", "2nd", "3rd", "Crew")
)
xtab
# Chi-square test
chisq_test(xtab)
# Compare the proportion of survived between groups
pairwise_prop_test(xtab)

# Test of independence using the data-frame interface
#####
df <- data.frame(
  gender = rep(c("M", "F"), each = 100),
  smoker = rep(c("yes", "no", "yes", "no"), times = c(30, 70, 60, 40))
)
# Positional columns
df %>% chisq_test(gender, smoker)
# Equivalent, using vars (keeps `correct` settable)
df %>% chisq_test(vars = c("gender", "smoker"), correct = FALSE)
```

Description

Compute Cliff's delta, a non-parametric effect size for the difference between two groups. It is the standardized version of the Mann-Whitney statistic and estimates the probability that a randomly drawn value from one group exceeds a randomly drawn value from the other, minus the reverse probability: $\delta = (\#\{x > y\} - \#\{x < y\}) / (n_1 n_2)$. It ranges from -1 to 1 and, unlike the rank-biserial r , is unaffected by ties beyond their contribution to the counts.

See the Datanovia tutorial [Wilcoxon Test in R](#) for a worked walkthrough.

Usage

```
cliff_delta(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  ci = FALSE,
  conf.level = 0.95,
  ci.type = "perc",
  nboot = 1000,
  ...,
  boot.parallel = getOption("boot.parallel", "no"),
  boot.ncpus = getOption("boot.ncpus", 1L)
)
```

Arguments

<code>data</code>	a data frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable and <code>group</code> is a factor with two or more levels.
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
<code>ci</code>	if TRUE, a percentile bootstrap confidence interval is computed and added as the columns <code>conf.low</code> and <code>conf.high</code> , as for cohens_d() and wilcox_effsize() .
<code>conf.level</code>	The level for the confidence interval.
<code>ci.type</code>	The type of confidence interval to use. Can be any of "norm", "basic", "perc", or "bca". Passed to <code>boot::boot.ci</code> .
<code>nboot</code>	The number of replications to use for bootstrap.
<code>...</code>	other arguments; accepted for interface compatibility with cohens_d() and wilcox_effsize() but not used (Cliff's delta has no test backend to forward

	them to). paired is rejected: the statistic is defined for two independent samples only.
boot.parallel	The type of parallel operation to be used when computing the bootstrap confidence interval. Allowed values are "no" (default), "multicore" and "snow". Passed to <code>boot()</code> . Defaults to <code>getOption("boot.parallel", "no")</code> , so it can also be set globally with <code>options(boot.parallel = "multicore")</code> . Only used when <code>ci = TRUE</code> .
boot.ncpus	Integer. The number of processes to be used in the parallel bootstrap. Defaults to <code>getOption("boot.ncpus", 1L)</code> . Note that <code>boot.parallel</code> has no effect unless <code>boot.ncpus > 1</code> . Only used when <code>ci = TRUE</code> .

Details

The magnitude thresholds are those of Romano et al. (2006): $\text{ldelta} < 0.147$ "negligible", < 0.33 "small", < 0.474 "medium", otherwise "large". Cliff's delta is algebraically identical to the rank-biserial correlation, so the point estimate equals `effectsize::rank_biserial()`.

Value

a tibble with one row per comparison and the columns `.y.`, `group1`, `group2`, `effsize` (Cliff's delta), `n1`, `n2` and `magnitude`; `conf.low` / `conf.high` are added when `ci = TRUE`.

References

- Cliff, N. (1993). Dominance statistics: Ordinal analyses to answer ordinal questions. *Psychological Bulletin*, 114(3), 494-509.
- Romano, J., Kromrey, J. D., Coraggio, J., & Skowronek, J. (2006). Appropriate statistics for ordinal level data. Annual meeting of the Florida Association of Institutional Research.

See Also

The Datanovia tutorial: [Wilcoxon Test in R](#).

Examples

```
# Two-samples Cliff's delta
ToothGrowth %>% cliff_delta(len ~ supp)

# Pairwise comparisons
ToothGrowth %>% cliff_delta(len ~ dose)

# Grouped data
ToothGrowth %>%
  dplyr::group_by(supp) %>%
  cliff_delta(len ~ dose)
```

cochran_qtest

*Cochran's Q Test***Description**

Performs the Cochran's Q test for unreplicated randomized block design experiments with a binary response variable and paired data. This test is analogue to the `friedman.test()` with 0,1 coded response. It's an extension of the McNemar Chi-squared test for comparing more than two paired proportions.

See the Datanovia tutorial [Cochran's Q Test in R](#) for a worked walkthrough.

Usage

```
cochran_qtest(data, formula)
```

Arguments

<code>data</code>	a data frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>a ~ b c</code> , where <code>a</code> is the outcome variable name; <code>b</code> is the within-subjects factor variables; and <code>c</code> (factor) is the column name containing individuals/subjects identifier. Should be unique per individual.

See Also

The Datanovia tutorial: [Cochran's Q Test in R](#).

Examples

```
# Generate a demo data
mydata <- data.frame(
  outcome = c(0,1,1,0,0,1,0,1,1,1,1,1,0,0,1,1,0,1,0,1,1,0,0,1,0,1,1,0,0,1),
  treatment = gl(3,1,30,labels=LETTERS[1:3]),
  participant = gl(10,3,labels=letters[1:10])
)
mydata$outcome <- factor(
  mydata$outcome, levels = c(1, 0),
  labels = c("success", "failure")
)
# Cross-tabulation
xtabs(~outcome + treatment, mydata)

# Compare the proportion of success between treatments
cochran_qtest(mydata, outcome ~ treatment|participant)

# pairwise comparisons between groups
pairwise_mcnemar_test(mydata, outcome ~ treatment|participant)
```

cohens_d

*Compute Cohen's d Measure of Effect Size***Description**

Compute the effect size for t-test. T-test conventional effect sizes, proposed by Cohen, are: 0.2 (small effect), 0.5 (moderate effect) and 0.8 (large effect).

Cohen's d is calculated as the difference between means or mean minus μ divided by the estimated standardized deviation.

For independent samples t-test, there are two possibilities implemented. If the t-test did not make a homogeneity of variance assumption, (the Welch test), the variance term will mirror the Welch test, otherwise a pooled estimate is used.

If a paired samples t-test was requested, then effect size desired is based on the standard deviation of the differences.

It can also return confidence intervals for the effect size, either by bootstrap (the default) or by an analytic, deterministic method (see `ci.method`).

See the Datanovia tutorial [Cohen's d Effect Size in R](#) for a worked walkthrough.

Usage

```
cohens_d(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  paired = FALSE,
  mu = 0,
  var.equal = FALSE,
  hedges.correction = FALSE,
  ci = FALSE,
  conf.level = 0.95,
  ci.type = "perc",
  nboot = 1000,
  boot.parallel = getOption("boot.parallel", "no"),
  boot.ncpus = getOption("boot.ncpus", 1L),
  id = NULL,
  ci.method = c("boot", "analytic")
)
```

Arguments

data	a data.frame containing the variables in the formula.
formula	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .

comparisons	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
ref.group	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
paired	a logical indicating whether you want a paired test.
mu	the theoretical mean (one-sample test) or the hypothesized difference in means (two-sample test). It is subtracted from the mean difference before standardizing, so a non-zero mu shifts the effect size accordingly. Default is 0.
var.equal	a logical variable indicating whether to treat the two variances as being equal. If TRUE then the pooled variance is used to estimate the variance otherwise the Welch (or Satterthwaite) approximation to the degrees of freedom is used. Used only for unpaired or independent samples test.
hedges.correction	logical indicating whether apply the Hedges correction by multiplying the usual value of Cohen's d by $(N-3)/(N-2.25)$ (for unpaired t-test) and by $(n1-2)/(n1-1.25)$ for paired t-test; where N is the total size of the two groups being compared ($N = n1 + n2$).
ci	If TRUE, returns confidence intervals by bootstrap. May be slow.
conf.level	The level for the confidence interval.
ci.type	The type of confidence interval to use. Can be any of "norm", "basic", "perc", or "bca". Passed to <code>boot::boot.ci</code> .
nboot	The number of replications to use for bootstrap.
boot.parallel	The type of parallel operation to be used when computing the bootstrap confidence interval. Allowed values are "no" (default), "multicore" and "snow". Passed to <code>boot()</code> . Defaults to <code>getOption("boot.parallel", "no")</code> , so it can also be set globally with <code>options(boot.parallel = "multicore")</code> . Only used when <code>ci = TRUE</code> .
boot.ncpus	Integer. The number of processes to be used in the parallel bootstrap. Defaults to <code>getOption("boot.ncpus", 1L)</code> . Note that <code>boot.parallel</code> has no effect unless <code>boot.ncpus > 1</code> . Only used when <code>ci = TRUE</code> .
id	(optional) character string with the name of the column holding the sample/subject identifier, used only for a paired test (<code>paired = TRUE</code>). When supplied, the two groups are matched by id (instead of by row order) before the paired d is computed, as in <code>t_test()</code> . Only complete pairs (subjects measured in both groups) are used.
ci.method	the method used to compute the confidence interval when <code>ci = TRUE</code> . Either "boot" (default) for a percentile bootstrap, or "analytic" for a deterministic interval obtained by inverting the noncentral <i>t</i> distribution (Steiger, 2004; Cumming & Finch, 2001) – the same machinery <code>anova_test(ci =)</code> uses for partial eta-squared. The analytic interval covers the one-sample, paired and two-sample (equal-variance and Welch) cases, is reproducible across runs (no seed), and

matches `effectsize::cohens_d(ci = )`. With `hedges.correction = TRUE` the estimate and both bounds are scaled by the documented $(N - 3)/(N - 2.25)$ approximation, so they are not identical to `effectsize::hedges_g(ci = )`, which applies the exact gamma-function correction at the design's degrees of freedom. The gap is proportional to the effect size and grows as the samples shrink. When the interval is not defined for a given input (for example a degenerate group), the bootstrap is used as a fallback. The default "boot" leaves the returned interval unchanged from previous versions.

Details

Quantification of the effect size magnitude is performed using the thresholds defined in Cohen (1992). The magnitude is assessed using the thresholds provided in (Cohen 1992), i.e. $|d| < 0.2$ "negligible", $|d| < 0.5$ "small", $|d| < 0.8$ "medium", otherwise "large".

Value

return a data frame with some of the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n, n1, n2`: Sample counts.
- `effsize`: estimate of the effect size (d value).
- `magnitude`: magnitude of effect size.
- `conf.low, conf.high`: lower and upper bound of the effect size confidence interval.

References

- Cohen, J. (1988). Statistical power analysis for the behavioral sciences (2nd ed.). New York:Academic Press.
- Cohen, J. (1992). A power primer. Psychological Bulletin, 112, 155-159.
- Hedges, Larry & Olkin, Ingram. (1985). Statistical Methods in Meta-Analysis. 10.2307/1164953.
- Navarro, Daniel. 2015. Learning Statistics with R: A Tutorial for Psychology Students and Other Beginners (Version 0.5).
- Steiger, J. H. (2004). Beyond the F test: Effect size confidence intervals and tests of close fit in the analysis of variance and contrast analysis. Psychological Methods, 9(2), 164-182.
- Cumming, G., & Finch, S. (2001). A primer on the understanding, use, and calculation of confidence intervals that are based on central and noncentral distributions. Educational and Psychological Measurement, 61(4), 532-574.

See Also

The Datanovia tutorial: [Cohen's d Effect Size in R](#).

Examples

```
# One-sample t test effect size
ToothGrowth %>% cohens_d(len ~ 1, mu = 0)

# Two indepedent samples t-test effect size
ToothGrowth %>% cohens_d(len ~ supp, var.equal = TRUE)

# Paired samples effect size
df <- data.frame(
  id = 1:5,
  pre = c(110, 122, 101, 120, 140),
  post = c(150, 160, 110, 140, 155)
)
df <- df %>% gather(key = "treatment", value = "value", -id)
head(df)

df %>% cohens_d(value ~ treatment, paired = TRUE)
```

conover_test

Conover's All-Pairs Rank Comparison Test

Description

Performs Conover's test (also known as the Conover-Iman test) for pairwise multiple comparisons of the ranked data, following a significant Kruskal-Wallis test. It is closely related to [dunn_test\(\)](#), but uses the pooled within-group rank variance and refers the test statistic to a t -distribution (with $N - k$ degrees of freedom) instead of the standard normal distribution. The Conover-Iman test is generally more powerful than Dunn's test, but should only be used as a post-hoc procedure when the Kruskal-Wallis test is itself significant (Conover, 1999).

If a reference group is specified (via `ref.group`), then each of the remaining group levels is compared only to the reference (control) group, and the p-value adjustment for multiple comparisons is computed over only these $k - 1$ comparisons (instead of all $k(k - 1)/2$ pairwise comparisons), exactly as for [dunn_test\(\)](#).

See the Datanovia tutorial [Kruskal-Wallis Test in R](#) for a worked walkthrough.

Usage

```
conover_test(
  data,
  formula,
  p.adjust.method = "holm",
  ref.group = NULL,
  detailed = FALSE
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference (control) group, and the p-value adjustment is computed over only these comparisons. Note that, like <code>dunn_test()</code> , <code>conover_test()</code> does not support <code>ref.group = "all"</code> .
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.

Details

The Conover-Iman pairwise statistic for comparing groups i and j is

$$t_{ij} = \frac{\bar{R}_i - \bar{R}_j}{\sqrt{S^2 \frac{N-1-H}{N-k} \left(\frac{1}{n_i} + \frac{1}{n_j} \right)}}$$

where $\bar{R}$ are the mean ranks, H is the (tie-corrected) Kruskal-Wallis statistic, N is the total sample size, k is the number of groups, and S^2 is the variance of the ranks ($S^2 = N(N+1)/12$ when there are no ties; otherwise $S^2 = \frac{1}{N-1} \left(\sum r^2 - \frac{N(N+1)^2}{4} \right)$). The statistic is referred to a t -distribution with $N - k$ degrees of freedom.

In the returned table each row is oriented with $i = \text{group2}$ and $j = \text{group1}$: estimate is $\bar{R}_{\text{group2}} - \bar{R}_{\text{group1}}$ and statistic carries its sign, the same convention as `dunn_test()`.

The results match `PMCMRplus::kwAllPairsConoverTest()`.

Value

return a data frame with some of the following columns:

- `.y`: the y (outcome) variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n1, n2`: Sample counts.
- `estimate`: mean ranks difference.
- `estimate1, estimate2`: show the mean rank values of the two groups, respectively.
- `statistic`: Test statistic (t-value) used to compute the p-value.
- `df`: degrees of freedom ($N - k$, the same for every comparison).
- `p`: p-value.

- p.adj: the adjusted p-value.
- method: the statistical test used to compare groups.
- p.adj.signif: the significance level of the adjusted p-values.

The **returned object has an attribute called args**, which is a list holding the test arguments.

References

Conover, W. J. (1999) Practical Nonparametric Statistics, 3rd edition. Wiley.

Conover, W. J. and Iman, R. L. (1979) On multiple-comparisons procedures. Technical Report LA-7677-MS, Los Alamos Scientific Laboratory.

See Also

[dunn_test](#), [kruskal_test](#) The Datanovia tutorial: [Kruskal-Wallis Test in R](#).

Examples

```
# Simple test
ToothGrowth %>% conover_test(len ~ dose)

# Comparison against a reference (control) group
# each group is compared to the reference; the p-value
# adjustment corrects for only these k - 1 comparisons
ToothGrowth %>% conover_test(len ~ dose, ref.group = "0.5")

# Grouped data
ToothGrowth %>%
  group_by(supp) %>%
  conover_test(len ~ dose)
```

convert_as_factor	<i>Factors</i>
-------------------	----------------

Description

Provides pipe-friendly functions to convert simultaneously multiple variables into a factor variable.

Helper functions are also available to set the reference level and the levels order.

Usage

```
convert_as_factor(data, ..., vars = NULL, make.valid.levels = FALSE)

set_ref_level(data, name, ref)

reorder_levels(data, name, order)
```


Arguments

<code>data</code>	a data frame
<code>...</code>	one unquoted expressions (or variable name) specifying the name of the variables you want to convert into factor. Alternative to the argument <code>vars</code> .
<code>vars</code>	a character vector specifying the variables to convert into factor.
<code>make.valid.levels</code>	logical. Default is FALSE. If TRUE, converts the variable to factor and add a leading character (x) if starting with a digit.
<code>name</code>	a factor variable name. Can be unquoted. For example, use <code>group</code> or <code>"group"</code> .
<code>ref</code>	the reference level.
<code>order</code>	a character vector specifying the order of the factor levels

Functions

- `convert_as_factor()`: Convert one or multiple variables into factor.
- `set_ref_level()`: Change a factor reference level or group.
- `reorder_levels()`: Change the order of a factor levels

Examples

```
# Create a demo data
df <- tibble(
  group = c("a", "a", "b", "b", "c", "c"),
  time = c("t1", "t2", "t1", "t2", "t1", "t2"),
  value = c(5, 6, 1, 3, 4, 5)
)
df
# Convert group and time into factor variable
result <- df %>% convert_as_factor(group, time)
result
# Show group levels
levels(result$group)

# Set c as the reference level (the first one)
result <- result %>%
  set_ref_level("group", ref = "c")
levels(result$group)

# Set the order of levels
result <- result %>%
  reorder_levels("group", order = c("b", "c", "a"))
levels(result$group)
```

cor_as_symbols

*Replace Correlation Coefficients by Symbols***Description**

Take a correlation matrix and replace the correlation coefficients by symbols according to the level of the correlation.

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

```
cor_as_symbols(
  x,
  cutpoints = c(0, 0.25, 0.5, 0.75, 1),
  symbols = c(" ", ".", "+", "*")
)
```

Arguments

x	a correlation matrix. Particularly, an object of class <code>cor_mat</code> .
cutpoints	numeric vector used for intervals. Default values are <code>c(0, 0.25, 0.5, 0.75, 1)</code> .
symbols	character vector, one shorter than cutpoints, used as correlation coefficient symbols. Default values are <code>c(" ", ".", "+", "*")</code> .

See Also

[cor_mat\(\)](#) The Datanovia tutorial: [Correlation Matrix in R](#).

Examples

```
# Compute correlation matrix
#::::::::::::::::::::::::::::::::::::
cor.mat <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat()

# Replace correlation coefficient by symbols
#::::::::::::::::::::::::::::::::::::
cor.mat %>%
  cor_as_symbols() %>%
  pull_lower_triangle()
```



```

cor.mat <- mydata %>% cor_mat()
cor.mat

# Collapse the correlation matrix into long format
# paired list data frame
long.format <- cor.mat %>% cor_gather()
long.format

# Spread a correlation data format
#::::::::::::::::::::::::::::::::::::::::::
# Spread the correlation coefficient value
long.format %>% cor_spread(value = "cor")
# Spread the p-value
long.format %>% cor_spread(value = "p")

```

cor_mark_significant *Add Significance Levels To a Correlation Matrix*

Description

Combines correlation coefficients and significance levels in a correlation matrix data.

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

```

cor_mark_significant(
  x,
  cutpoints = c(0, 1e-04, 0.001, 0.01, 0.05, 1),
  symbols = c("****", "***", "**", "*", "")
)

```

Arguments

x	an object of class <code>cor_mat()</code> .
cutpoints	numeric vector used for intervals.
symbols	character vector, one shorter than cutpoints, used as significance symbols.

Value

a data frame containing the lower triangular part of the correlation matrix marked by significance symbols.

See Also

The Datanovia tutorial: [Correlation Matrix in R](#).

Examples

```
mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat() %>%
  cor_mark_significant()
```

cor_mat

Compute Correlation Matrix with P-values

Description

Compute correlation matrix with p-values. Numeric columns in the data are detected and automatically selected for the analysis. You can also specify variables of interest to be used in the correlation analysis.

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

```
cor_mat(
  data,
  ...,
  vars = NULL,
  method = "pearson",
  alternative = "two.sided",
  conf.level = 0.95
)

cor_pmat(
  data,
  ...,
  vars = NULL,
  method = "pearson",
  alternative = "two.sided",
  conf.level = 0.95
)

cor_get_pval(x)
```

Arguments

data	a data.frame containing the variables.
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest.
vars	a character vector containing the variable names of interest.
method	a character string indicating which correlation coefficient is to be used for the test. One of "pearson", "kendall", or "spearman", can be abbreviated.

method	Character, the visualization method of correlation matrix to be used. Currently, it supports seven methods, named 'circle' (default), 'square', 'ellipse', 'number', 'pie', 'shade' and 'color'. See examples for details. The areas of circles or squares show the absolute value of corresponding correlation coefficients. Method 'pie' and 'shade' came from Michael Friendly's job (with some adjustment about the shade added on), and 'ellipse' came from D.J. Murdoch and E.D. Chow's job, see in section References.
type	Character, 'full' (default), 'upper' or 'lower', display full matrix, lower triangular or upper triangular matrix.
palette	character vector containing the color palette.
p.mat	matrix of p-value corresponding to the correlation matrix.
significant.level	significant level, if the p-value is bigger than significant.level, then the corresponding correlation coefficient is regarded as insignificant.
insignificant	character, specialized insignificant correlation coefficients, "cross" (default), "blank". If "blank", wipe away the corresponding glyphs; if "cross", add crosses (X) on corresponding glyphs.
label	logical value. If TRUE, shows the correlation coefficient labels.
font.label	a list with one or more of the following elements: size (e.g., 1), color (e.g., "black") and style (e.g., "bold"). Used to customize the correlation coefficient labels. For example font.label = list(size = 1, color = "black", style = "bold").
...	additional options not listed (i.e. "tl.cex") here to pass to corrplot.

See Also

[cor_as_symbols\(\)](#) The Datanovia tutorial: [Correlation Matrix in R](#).

Examples

```
# Compute correlation matrix
#::::::::::::::::::::::::::::::::::::
cor.mat <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat()

# Visualize correlation matrix
#::::::::::::::::::::::::::::::::::::
# Full correlation matrix,
# insignificant correlations are marked by crosses
cor.mat %>% cor_plot()

# Reorder by correlation coefficient
# pull lower triangle and visualize
cor.lower.tri <- cor.mat %>%
  cor_reorder() %>%
  pull_lower_triangle()
cor.lower.tri %>% cor_plot()
```



```

# Change visualization methods
#::::::::::::::::::::::::::::::::::::
cor.lower.tri %>%
  cor_plot(method = "pie")

cor.lower.tri %>%
  cor_plot(method = "color")

cor.lower.tri %>%
  cor_plot(method = "number")

# Show the correlation coefficient: label = TRUE
# Blank the insignificant correlation
#::::::::::::::::::::::::::::::::::::
cor.lower.tri %>%
  cor_plot(
    method = "color",
    label = TRUE,
    insignificant = "blank"
  )

# Change the color palettes
#::::::::::::::::::::::::::::::::::::

# Using custom color palette
# Require ggpubr: install.packages("ggpubr")
if(require("ggpubr")){
  my.palette <- get_palette(c("red", "white", "blue"), 200)
  cor.lower.tri %>%
    cor_plot(palette = my.palette)
}

# Using RcolorBrewer color palette
if(require("ggpubr")){
  my.palette <- get_palette("PuOr", 200)
  cor.lower.tri %>%
    cor_plot(palette = my.palette)
}

```

cor_reorder

Reorder Correlation Matrix

Description

reorder correlation matrix, according to the coefficients, using the hierarchical clustering method.

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

cor_reorder(x)

Arguments

x a correlation matrix. Particularly, an object of class cor_mat.

Value

a data frame

See Also

[cor_mat\(\)](#), [cor_gather\(\)](#), [cor_spread\(\)](#) The Datanovia tutorial: [Correlation Matrix in R](#).

Examples

```
# Compute correlation matrix
#::::::::::::::::::::::::::::::::::::::::::::
cor.mat <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat()

# Reorder by correlation and get p-values
#::::::::::::::::::::::::::::::::::::::::::::
# Reorder
cor.mat %>%
  cor_reorder()
# Get p-values of the reordered cor_mat
cor.mat %>%
  cor_reorder() %>%
  cor_get_pval()
```

cor_select	<i>Subset Correlation Matrix</i>
------------	----------------------------------

Description

See the Datanovia tutorial [Correlation Matrix in R](#) for a worked walkthrough.

Usage

cor_select(x, ..., vars = NULL)

Arguments

x	a correlation matrix. Particularly, an object of class <code>cor_mat</code> .
...	One or more unquoted expressions (or variable names) separated by commas. Used to select variables of interest.
vars	a character vector containing the variable names of interest.

Value

a data frame

See Also

`cor_mat()`, `pull_triangle()`, `replace_triangle()`. The Datanovia tutorial: [Correlation Matrix in R](#).

Examples

```
# Compute correlation matrix
#::::::::::::::::::::::::::::::::::::::::::::
cor.mat <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat()

# Subsetting correlation matrix
#::::::::::::::::::::::::::::::::::::::::::::

# Select some variables of interest
cor.mat %>%
  cor_select(mpg, drat, wt)

# Remove variables
cor.mat %>%
  cor_select(-mpg, -wt)
```

cor_test

Correlation Test

Description

Provides a pipe-friendly framework to perform correlation test between paired samples, using Pearson, Kendall or Spearman method. Wrapper around the function `cor.test()`.

Can also performs multiple pairwise correlation analyses between more than two variables or between two different vectors of variables. Using this function, you can also compute, for example, the correlation between one variable vs many.

See the Datanovia tutorial [Correlation Test in R](#) for a worked walkthrough.

Usage

```
cor_test(
  data,
  ...,
  vars = NULL,
  vars2 = NULL,
  alternative = "two.sided",
  method = "pearson",
  conf.level = 0.95,
  use = "pairwise.complete.obs"
)
```

Arguments

<code>data</code>	a data.frame containing the variables.
<code>...</code>	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest. Alternative to the argument <code>vars</code> .
<code>vars</code>	optional character vector containing variable names for correlation analysis. Ignored when dot <code>vars</code> are specified. • If <code>vars</code> is <code>NULL</code>, multiple pairwise correlation tests is performed between all variables in the data. • If <code>vars</code> contain only one variable, a pairwise correlation analysis is performed between the specified variable vs either all the remaining numeric variables in the data or variables in <code>vars2</code> (if specified). • If <code>vars</code> contain two or more variables: i) if <code>vars2</code> is not specified, a pairwise correlation analysis is performed between all possible combinations of variables. ii) if <code>vars2</code> is specified, each element in <code>vars</code> is tested against all elements in <code>vars2</code> . Accept unquoted variable names: <code>c(var1, var2)</code>.
<code>vars2</code>	optional character vector. If specified, each element in <code>vars</code> is tested against all elements in <code>vars2</code> . Accept unquoted variable names: <code>c(var1, var2)</code> .
<code>alternative</code>	indicates the alternative hypothesis and must be one of "two.sided", "greater" or "less". You can specify just the initial letter. "greater" corresponds to positive association, "less" to negative association.
<code>method</code>	a character string indicating which correlation coefficient is to be used for the test. One of "pearson", "kendall", or "spearman", can be abbreviated.
<code>conf.level</code>	confidence level for the returned confidence interval. Currently only used for the Pearson product moment correlation coefficient if there are at least 4 complete pairs of observations.
<code>use</code>	an optional character string giving a method for computing covariances in the presence of missing values. This must be (an abbreviation of) one of the strings "everything", "all.obs", "complete.obs", "na.or.complete", or "pairwise.complete.obs".

Value

return a data frame with the following columns:

- var1, var2: the variables used in the correlation test.
- cor: the correlation coefficient.
- statistic: Test statistic used to compute the p-value.
- df: the degrees of freedom (Pearson method only).
- p: p-value.
- conf.low, conf.high: Lower and upper bounds on a confidence interval.
- method: the method used to compute the statistic.

Functions

- cor_test(): correlation test between two or more variables.

Note

cor_test() does not support weighted correlations: it wraps [cor.test\(\)](#), which has no weights argument. Passing weights = therefore raises an error rather than silently returning the unweighted result. For a weighted Pearson correlation use base R, e.g. `stats::cov.wt(data[, c("x", "y")], wt = data$w, cor = TRUE)$cor`.

See Also

[cor_mat\(\)](#), [as_cor_mat\(\)](#), [rstatix-programming](#) (using variable names held in strings) The Datanovia tutorial: [Correlation Test in R](#).

Examples

```
# Correlation between the specified variable vs
# the remaining numeric variables in the data
#::::::::::::::::::::::::::::::::::::::::::::
mtcars %>% cor_test(mpg)

# Correlation test between two variables
#::::::::::::::::::::::::::::::::::::::::::::
mtcars %>% cor_test(wt, mpg)

# Pairwise correlation between multiple variables
#::::::::::::::::::::::::::::::::::::::::::::
mtcars %>% cor_test(wt, mpg, disp)

# Grouped data
#::::::::::::::::::::::::::::::::::::::::::::
iris %>%
  group_by(Species) %>%
  cor_test(Sepal.Width, Sepal.Length)

# Multiple correlation test
```



```
# Convert into a data frame of cases
#####
df <- counts_to_cases(xtab)
head(df)
```

cramer_v

Compute Cramer's V

Description

Compute Cramer's V, which measures the strength of the association between categorical variables. See the Datanovia tutorial [Chi-Square Test of Independence in R](#) for a worked walkthrough.

Usage

```
cramer_v(x, y = NULL, correct = FALSE, ..., ci = FALSE, conf.level = 0.95)
```

Arguments

x	a numeric vector or matrix. x and y can also both be factors.
y	a numeric vector; ignored if x is a matrix. If x is a factor, y should be a factor of the same length.
correct	logical. If TRUE, Yates' continuity correction is applied when computing the chi-square statistic, which only affects 2x2 tables. Default is FALSE. Yates' correction improves the chi-square approximation to the <i>null</i> distribution of the test statistic, so it belongs to the <i>test</i> (see chisq.test()) rather than to an effect size: it shrinks the statistic and therefore biases Cramer's V downward. Set <code>correct = TRUE</code> to reproduce the value returned by earlier versions of <i>rstatix</i> , which applied the correction by default.
...	other arguments passed to the function chisq.test() .
ci	logical. If TRUE, a confidence interval for Cramer's V is added to the result and a one-row data frame is returned instead of a single value. Default is FALSE.
conf.level	The level of the confidence interval. Default is 0.95. Only used when <code>ci = TRUE</code> .

Details

Cramer's V is $V = \sqrt{\chi^2 / (N(k - 1))}$ (Cramer, 1946), where χ^2 is the Pearson chi-square statistic, N the total count and k the smaller of the two table dimensions.

The confidence interval is obtained by inverting the noncentral chi-square distribution (Smithson, 2003; Steiger, 2004): the noncentrality parameters λ whose distributions place the observed chi-square statistic at the $1 - \alpha/2$ and $\alpha/2$ quantiles are found by root finding, and each is converted with $V = \sqrt{\lambda / (N(k - 1))}$. The interval is computed from the same chi-square statistic as the point estimate, so `correct = TRUE` shifts both.

The bounds are clipped to $[0, 1]$, the range of Cramer's V. They are NA, with a warning, when the statistic or its degrees of freedom are undefined – for instance when the table has an empty row or

column, or when `simulate.p.value = TRUE` is passed on to `chisq.test()`, which then reports no degrees of freedom.

The interval usually brackets the reported `effsize`, but it does not in the near-independence corner: when the observed chi-square falls below the $\alpha/2$ quantile of its central distribution, no noncentrality is consistent with the data at that quantile, both bounds collapse to $[0, 0]$, and the (necessarily positive) point estimate lies above them. This is a property of the noncentral inversion rather than of this implementation, and it only arises for effect sizes indistinguishable from zero.

At the default `correct = FALSE`, the results match `effectsize::cramers_v(adjust = FALSE, ci = , alternative = "two.sided")` away from the near-independence corner above (where numerical inversions differ), and `DescTools::CramerV(conf.level = )` to about four decimals (its inversion uses a looser tolerance), except at a chi-square of exactly zero, where `DescTools` returns NA bounds and this function returns the collapsed $[0, 0]$ interval. Neither package applies Yates' continuity correction, so `correct = TRUE` values have no counterpart there (`DescTools`'s own `correct` argument selects the Bergsma bias correction, a different adjustment).

Value

By default, a single numeric value: Cramer's V.

When `ci = TRUE`, a one-row tibble with the columns `effsize` (Cramer's V), `conf.low` and `conf.high` – the same confidence interval columns that `anova_test(ci = )` returns.

References

Cramer, H. (1946). *Mathematical Methods of Statistics*. Princeton University Press.

Smithson, M. (2003). *Confidence Intervals*. Sage Publications.

Steiger, J. H. (2004). Beyond the F test: Effect size confidence intervals and tests of close fit in the analysis of variance and contrast analysis. *Psychological Methods*, 9, 164-182.

See Also

The Datanovia tutorial: [Chi-Square Test of Independence in R](#).

Examples

```
# Data preparation
df <- as.table(rbind(c(762, 327, 468), c(484, 239, 477)))
dimnames(df) <- list(
  gender = c("F", "M"),
  party = c("Democrat", "Independent", "Republican")
)
df
# Compute cramer's V
cramer_v(df)

# Add a confidence interval
cramer_v(df, ci = TRUE)

# Yates' continuity correction only affects 2x2 tables. It belongs to the
# test, not to the effect size, so it is off by default.
```



```
tab <- as.table(rbind(c(20, 30), c(35, 15)))
cramer_v(tab)
cramer_v(tab, correct = TRUE)
```

df_arrange	<i>Arrange Rows by Column Values</i>
------------	--------------------------------------

Description

Order the rows of a data frame by values of specified columns. Wrapper around the [arrange\(\)](#) function. Supports standard and non standard evaluation.

Usage

```
df_arrange(data, ..., vars = NULL, .by_group = FALSE)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest. Use desc() to sort a variable in descending order.
vars	a character vector containing the variable names of interest.
.by_group	If TRUE, will sort first by grouping variable. Applies to grouped data frames only.

Value

a data frame

Examples

```
df <- head(ToothGrowth)
df

# Select column using standard evaluation
df %>% df_arrange(vars = c("dose", "len"))

# Select column using non-standard evaluation
df %>% df_arrange(dose, desc(len))
```

df_get_var_names	<i>Get User Specified Variable Names</i>
------------------	--

Description

Returns user specified variable names. Supports standard and non standard evaluation.

Usage

```
df_get_var_names(data, ..., vars = NULL)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest.
vars	a character vector containing the variable names of interest.

Value

a character vector

Examples

```
# Non standard evaluation
ToothGrowth %>%
  df_get_var_names(dose, len)

# Standard evaluation
ToothGrowth %>%
  df_get_var_names(vars = c("len", "dose"))
```

df_group_by	<i>Group a Data Frame by One or more Variables</i>
-------------	--

Description

Group a data frame by one or more variables. Supports standard and non standard evaluation.

Usage

```
df_group_by(data, ..., vars = NULL)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest.
vars	a character vector containing the variable names of interest.

Examples

```
# Non standard evaluation
by_dose <- head(ToothGrowth) %>%
  df_group_by(dose)
by_dose

# Standard evaluation
head(ToothGrowth) %>%
  df_group_by(vars = c("dose", "supp"))
```

df_label_both	<i>Functions to Label Data Frames by Grouping Variables</i>
---------------	---

Description

Functions to label data frame rows by one or multiple grouping variables.

Usage

```
df_label_both(data, ..., vars = NULL, label_col = "label", sep = c(", ", ":", ""))

df_label_value(data, ..., vars = NULL, label_col = "label", sep = ", ")
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used as grouping variables.
vars	a character vector containing the grouping variables of interest.
label_col	column to hold the label of the data subsets. Default column name is "label".
sep	String separating labelling variables and values. Should be of length 2 in the function df_label_both(). 1) One sep is used to separate groups, for example ','; 2) The other sep between group name and levels; for example ':'.

Value

a modified data frame with a column containing row labels.

Functions

- `df_label_both()`: Displays both the variable name and the factor value.
- `df_label_value()`: Displays only the value of a factor.

Examples

```
# Data preparation
df <- head(ToothGrowth)

# Labelling: Non standard evaluation
df %>%
  df_label_both(dose, supp)

# Standard evaluation
df %>%
  df_label_both(dose, supp)

# Nesting the data then label each subset by groups
ToothGrowth %>%
  df_nest_by(dose, supp) %>%
  df_label_both(supp, dose)
```

df_nest_by

Nest a Tibble By Groups

Description

Nest a tibble data frame using grouping specification. Supports standard and non standard evaluation.

Usage

```
df_nest_by(data, ..., vars = NULL)
```

Arguments

<code>data</code>	a data frame
<code>...</code>	One or more unquoted expressions (or variable names) separated by commas. Used as grouping variables.
<code>vars</code>	a character vector containing the grouping variables of interest.

Value

A tibble with one row per unique combination of the grouping variables. The first columns are the grouping variables, followed by a list column of tibbles with matching rows of the remaining columns.

Examples

```
# Non standard evaluation
ToothGrowth %>%
  df_nest_by(dose, supp)

# Standard evaluation
ToothGrowth %>%
  df_nest_by(vars = c("dose", "supp"))
```

df_select

*Select Columns in a Data Frame***Description**

A wrapper around the [select\(\)](#) function for selection data frame columns. Supports standard and non standard evaluations. Usefull to easily program with dplyr

Usage

```
df_select(data, ..., vars = NULL)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest.
vars	a character vector containing the variable names of interest.

Value

a data frame

Examples

```
df <- head(ToothGrowth)
df

# Select column using standard evaluation
df %>% df_select(vars = c("dose", "len"))

# Select column using non-standard evaluation
df %>% df_select(dose, len)
```

df_split_by

*Split a Data Frame into Subset***Description**

Split a data frame by groups into subsets or data panel. Very similar to the function `df_nest_by()`. The only difference is that, it adds label to each data subset. Labels are the combination of the grouping variable levels. The column holding labels are named "label".

Usage

```
df_split_by(
  data,
  ...,
  vars = NULL,
  label_col = "label",
  labeller = df_label_both,
  sep = c(", ", ":",)
)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used as grouping variables.
vars	a character vector containing the grouping variables of interest.
label_col	column to hold the label of the data subsets. Default column name is "label".
labeller	A function that takes a data frame, the grouping variables, label_col and label_sep arguments, and add labels into the data frame. Example of possible values are: <code>df_label_both()</code> and <code>df_label_value()</code> .
sep	String separating labelling variables and values. Should be of length 2 in the function <code>df_label_both()</code> . 1) One sep is used to separate groups, for example ','; 2) The other sep between group name and levels; for example ':'.

Value

A tibble with one row per unique combination of the grouping variables. The first columns are the grouping variables, followed by a list column of tibbles with matching rows of the remaining columns, and a column named label, containing labels.

Examples

```
# Split a data frame
# .....
# Create a grouped data
res <- ToothGrowth %>%
```

```

    df_split_by(dose, supp)
  res

  # Show subsets
  res$data

  # Add panel/subset labels
  res <- ToothGrowth %>%
    df_split_by(dose, supp)
  res

```

df_unite	<i>Unite Multiple Columns into One</i>
----------	--

Description

Paste together multiple columns into one. Wrapper around `unite()` that supports standard and non standard evaluation.

Usage

```

df_unite(data, col, ..., vars = NULL, sep = "_", remove = TRUE, na.rm = FALSE)

df_unite_factors(
  data,
  col,
  ...,
  vars = NULL,
  sep = "_",
  remove = TRUE,
  na.rm = FALSE
)

```

Arguments

data	a data frame
col	the name of the new column as a string or a symbol.
...	a selection of columns. One or more unquoted expressions (or variable names) separated by commas.
vars	a character vector containing the column names of interest.
sep	Separator to use between values.
remove	If TRUE, remove input columns from output data frame.
na.rm	If TRUE, missing values will be removed prior to uniting each value.

Functions

- `df_unite()`: Unite multiple columns into one.
- `df_unite_factors()`: Unite factor columns. First, order factors levels then merge them into one column. The output column is a factor.

Examples

```
# Non standard evaluation
head(ToothGrowth) %>%
  df_unite(col = "dose_supp", dose, supp)

# Standard evaluation
head(ToothGrowth) %>%
  df_unite(col = "dose_supp", vars = c("dose", "supp"))
```

doo

Alternative to dplyr::do for Doing Anything

Description

Provides a flexible alternative to the `dplyr::do()` function. Technically it uses `nest()` + `mutate()` + `map()` to apply arbitrary computation to a grouped data frame.

The output is a data frame. If the applied function returns a data frame, then the output will be automatically unnested. Otherwise, the output includes the grouping variables and a column named `".results."` (by default), which is a "list-columns" containing the results for group combinations.

Usage

```
doo(data, .f, ..., result = ".results.")
```

Arguments

<code>data</code>	a (grouped) data frame
<code>.f</code>	A function, formula, or atomic vector. For example <code>~t.test(len ~ supp, data = .)</code> .
<code>...</code>	Additional arguments passed on to <code>.f</code>
<code>result</code>	the column name to hold the results. Default is <code>".results."</code> .

Value

a data frame

Usage

```
dunnett_test(
  data,
  formula,
  ref.group = NULL,
  conf.level = 0.95,
  detailed = FALSE
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>ref.group</code>	a character string specifying the reference (control) group. Each remaining group level is compared against this group. If <code>NULL</code> (default), the first level of the grouping variable is used as the control.
<code>conf.level</code>	confidence level of the (simultaneous) confidence intervals.
<code>detailed</code>	logical value. Default is <code>FALSE</code> . If <code>TRUE</code> , a detailed result is shown.

Value

a data frame with some of the following columns:

- `.y.`: the outcome variable used in the test.
- `group1, group2`: the compared groups; `group1` is the control (reference) and `group2` is the treatment, consistent with the `ref.group` convention of `t_test()`/`wilcox_test()`/`dunn_test()`/`emmeans_test()`.
- `n1, n2`: sample sizes of the control and treatment groups.
- `estimate`: the estimated mean difference `group1 - group2` (control minus treatment).
- `conf.low, conf.high`: simultaneous confidence interval for the difference.
- `statistic`: the t-statistic.
- `df`: degrees of freedom.
- `p.adj`: the Dunnett-adjusted p-value.
- `method`: the statistical test used.
- `p.adj.signif`: the significance level of the adjusted p-value.

The `estimate`, `confidence interval`, `se` and `method` columns are returned only when `detailed = TRUE`.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

References

Dunnett, C. W. (1955) A multiple comparison procedure for comparing several treatments with a control. *Journal of the American Statistical Association*, 50, 1096-1121.

See Also

`tukey_hsd()`, `games_howell_test()`, `emmeans_test()` The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
if (requireNamespace("emmeans", quietly = TRUE)) {

  # Compare each dose to the control dose ("0.5")
  ToothGrowth %>% dunnett_test(len ~ dose)

  # Detailed output (estimate + simultaneous confidence interval)
  ToothGrowth %>% dunnett_test(len ~ dose, detailed = TRUE)

  # Grouped data
  ToothGrowth %>%
    group_by(supp) %>%
    dunnett_test(len ~ dose)

}
```

dunn_test

*Dunn's Test of Multiple Comparisons***Description**

Performs Dunn's test for pairwise multiple comparisons of the ranked data. The mean rank of the different groups is compared. Used for post-hoc test following Kruskal-Wallis test.

The default of the `rstatix::dunn_test()` function is to perform a two-sided Dunn test like the well known commercial softwares, such as SPSS and GraphPad. This is not the case for some other R packages (`dunn.test` and `jamovi`), where the default is to perform one-sided test. This discrepancy is documented at <https://github.com/kassambara/rstatix/issues/50>.

If a reference group is specified (via `ref.group`), then each of the remaining group levels is compared only to the reference (control) group, and the p-value adjustment for multiple comparisons is computed over only these $k - 1$ comparisons (instead of all $k(k - 1)/2$ pairwise comparisons). Note that this affects the adjusted p-values: it is not equivalent to filtering the full pairwise result afterwards, which would still adjust over all pairs.

See the Datanovia tutorial [Kruskal-Wallis Test in R](#) for a worked walkthrough.

Usage

```
dunn_test(
  data,
  formula,
  p.adjust.method = "holm",
  ref.group = NULL,
  detailed = FALSE,
```

```
    effect.size = FALSE
  )
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference (control) group, and the p-value adjustment is computed over only these comparisons. Note that, unlike <code>t_test()</code> and <code>wilcox_test()</code> , <code>dunn_test()</code> does not support <code>ref.group = "all"</code> .
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>effect.size</code>	logical. Default is FALSE. If TRUE, an <code>r</code> column is added, the effect size $r = Z / \sqrt{N}$ where Z is Dunn's z-statistic and N is the total sample size (the whole-sample rank variance Dunn's z is standardised on, not the pairwise $n1 + n2$). No magnitude label is attached: there is no threshold set calibrated for it.

Details

DunnTest performs the post hoc pairwise multiple comparisons procedure appropriate to follow up a Kruskal-Wallis test, which is a non-parametric analog of the one-way ANOVA. The Wilcoxon rank sum test, itself a non-parametric analog of the unpaired t-test, is possibly intuitive, but inappropriate as a post hoc pairwise test, because (1) it fails to retain the dependent ranking that produced the Kruskal-Wallis test statistic, and (2) it does not incorporate the pooled variance estimate implied by the null hypothesis of the Kruskal-Wallis test.

Value

return a data frame with some of the following columns:

- `.y.`: the `y` (outcome) variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n1, n2`: Sample counts.
- `estimate`: mean ranks difference.
- `estimate1, estimate2`: show the mean rank values of the two groups, respectively.
- `statistic`: Test statistic (z-value) used to compute the p-value.
- `p`: p-value.
- `p.adj`: the adjusted p-value.

- method: the statistical test used to compare groups.
- p.signif, p.adj.signif: the significance level of p-values and adjusted p-values, respectively.

The **returned object has an attribute called args**, which is a list holding the test arguments.

References

Dunn, O. J. (1964) Multiple comparisons using rank sums *Technometrics*, 6(3):241-252.

See Also

The Datanovia tutorial: [Kruskal-Wallis Test in R](#).

Examples

```
# Simple test
ToothGrowth %>% dunn_test(len ~ dose)

# Comparison against a reference (control) group
# each group is compared to the reference; the p-value
# adjustment corrects for only these k - 1 comparisons
ToothGrowth %>% dunn_test(len ~ dose, ref.group = "0.5")

# Grouped data
ToothGrowth %>%
  group_by(supp) %>%
  dunn_test(len ~ dose)
```

emmeans_test

Pairwise Comparisons of Estimated Marginal Means

Description

Performs pairwise comparisons between groups using the estimated marginal means. Pipe-friendly wrapper around the functions `emmeans()` + `contrast()` from the `emmeans` package, which need to be installed before using this function. This function is useful for performing post-hoc analyses following ANOVA/ANCOVA tests.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
emmeans_test(
  data,
  formula,
  covariate = NULL,
  ref.group = NULL,
  comparisons = NULL,
```

```

    p.adjust.method = "bonferroni",
    conf.level = 0.95,
    model = NULL,
    detailed = FALSE
)

get_emmeans(emmeans.test)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>covariate</code>	(optional) covariate names (for ANCOVA)
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>conf.level</code>	confidence level of the interval.
<code>model</code>	a fitted-model object such as the result of a call to <code>lm()</code> , <code>stats::aov()</code> (including a within-subject <code>Error()</code> term) or <code>nlme::lme()</code> , on which the estimated marginal means are computed. Supplying <code>model</code> lets you (i) average over factors that are in the model but not in <code>formula</code> (e.g. for a factorial design, fit <code>lm(y ~ a * b)</code> and compare <code>a</code> with <code>formula = y ~ a</code> , averaging over <code>b</code>), and (ii) run pairwise comparisons for repeated-measures / mixed designs (pass a within-subject model). When <code>model = NULL</code> (default), a simple <code>lm()</code> is fitted from <code>formula</code> (plus the grouping and covariate variables). See examples.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>emmeans.test</code>	an object of class <code>emmeans_test</code> .

Value

return a data frame with some the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.

- `statistic`: Test statistic (t.ratio) used to compute the p-value.
- `df`: degrees of freedom.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the statistical test used to compare groups.
- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: estimate of the effect size, that is the difference between the two emmeans (estimated marginal means).
- `conf.low`, `conf.high`: Lower and upper bound on a confidence interval of the estimate.

The **returned object has an attribute called `args`**, which is a list holding the test arguments. It has also an attribute named "emmeans", a data frame containing the groups emmeans.

Functions

- `get_emmeans()`: returns the estimated marginal means from an object of class `emmeans_test`

See Also

The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
if (requireNamespace("emmeans", quietly = TRUE)) {

  # Data preparation
  df <- ToothGrowth
  df$dose <- as.factor(df$dose)

  # Pairwise comparisons
  res <- df %>%
    group_by(supp) %>%
    emmeans_test(len ~ dose, p.adjust.method = "bonferroni")
  res

  # Display estimated marginal means
  attr(res, "emmeans")

  # Show details
  df %>%
    group_by(supp) %>%
    emmeans_test(len ~ dose, p.adjust.method = "bonferroni", detailed = TRUE)

  # Marginal means averaged over another factor (e.g. a 2x3 design).
  # Fit the full model and pass it with `model =` so that the estimated
  # marginal means for `dose` are averaged over `supp` (instead of fitting
  # `len ~ dose` alone, which would ignore `supp`):
  model <- lm(len ~ supp * dose, data = df)
  df %>% emmeans_test(len ~ dose, model = model)
```

```

# Repeated-measures / mixed designs: pass a fitted within-subject model
# (e.g. stats::aov() with an Error() term, or nlme::lme()) with `model`=:

set.seed(123)
d <- data.frame(
  id    = factor(rep(1:10, 3)),
  time  = factor(rep(c("t1", "t2", "t3"), each = 10)),
  score = rnorm(30)
)
rm_model <- stats::aov(score ~ time + Error(id / time), data = d)
d %>% emmeans_test(score ~ time, model = rm_model)

}

```

eta_squared

*Effect Size for ANOVA***Description**

Compute eta-squared and partial eta-squared for all terms in an ANOVA model.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```

eta_squared(model, ci = NULL)

partial_eta_squared(model, ci = NULL)

```

Arguments

<code>model</code>	an object of class <code>aov</code> or <code>anova</code> .
<code>ci</code>	confidence level for a confidence interval on the effect size. If a number between 0 and 1 (e.g. 0.95), the function returns a tibble with one row per model term and the columns <code>Effect</code> , <code>effsize</code> , <code>conf.low</code> and <code>conf.high</code> instead of the bare named vector. The interval is computed in base R by inverting the noncentral F distribution (Steiger, 2004), and matches <code>effectsize::eta_squared(ci = , alternative = "two.sided")</code> — <code>partial = FALSE</code> for <code>eta_squared()</code> and <code>partial = TRUE</code> for <code>partial_eta_squared()</code> — to about four decimals for the non-partial bounds of a small pseudo-F (that function's inversion uses a looser tolerance), and more closely everywhere else. Default is <code>NULL</code> (no interval; the bare named vector is returned, unchanged).

Value

a named numeric vector of effect sizes, one per model term; or, when `ci` is a confidence level, a tibble with the columns `Effect`, `effsize`, `conf.low` and `conf.high`.

Functions

- `eta_squared()`: compute eta squared
- `partial_eta_squared()`: compute partial eta squared.

References

Steiger, J. H. (2004). Beyond the F test: Effect size confidence intervals and tests of close fit in the analysis of variance and contrast analysis. *Psychological Methods*, 9, 164-182.

See Also

The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Data preparation
df <- ToothGrowth
df$dose <- as.factor(df$dose)

# Compute ANOVA
res.aov <- aov(len ~ supp*dose, data = df)
summary(res.aov)

# Effect size
eta_squared(res.aov)
partial_eta_squared(res.aov)

# Effect size with confidence interval
eta_squared(res.aov, ci = 0.95)
partial_eta_squared(res.aov, ci = 0.95)
```

factorial_design

Build Factorial Designs for ANOVA

Description

Provides helper functions to build factorial design for easily computing ANOVA using the [Anova\(\)](#) function. This might be very useful for repeated measures ANOVA, which is hard to set up with the `car` package.

Usage

```
factorial_design(data, dv, wid, between, within, covariate)
```



```

df$id <- rep(1:10, 6) # Add individuals id
head(df)
# Build factorial designs
design <- factorial_design(df, dv = len, wid = id, within = c(supp, dose))
design
# Easily perform repeated measures ANOVA using the car package
res.anova <- Anova(design$model, idata = design$idata, idesign = design$idesign, type = 3)
summary(res.anova, multivariate = FALSE)

# Independent measures designs
#::::::::::::::::::::::::::::::::::::
# Build factorial designs
df$id <- 1:nrow(df)
design <- factorial_design(df, dv = len, wid = id, between = c(supp, dose))
design
# Perform ANOVA
Anova(design$model, type = 3)

```

fisher_test

Fisher's Exact Test for Count Data

Description

Performs Fisher's exact test for testing the null of independence of rows and columns in a contingency table.

Wrappers around the R base function `fisher.test()` but have the advantage of performing pairwise and row-wise fisher tests, the post-hoc tests following a significant chi-square test of homogeneity for 2xc and rx2 contingency tables.

See the Datanovia tutorial [Fisher's Exact Test in R](#) for a worked walkthrough.

Usage

```

fisher_test(
  xtab,
  workspace = 2e+05,
  alternative = "two.sided",
  conf.int = TRUE,
  conf.level = 0.95,
  simulate.p.value = FALSE,
  B = 2000,
  detailed = FALSE,
  ...
)

pairwise_fisher_test(xtab, p.adjust.method = "holm", detailed = FALSE, ...)

row_wise_fisher_test(xtab, p.adjust.method = "holm", detailed = FALSE, ...)

```

Arguments

<code>xtab</code>	a contingency table in a matrix form.
<code>workspace</code>	an integer specifying the size of the workspace used in the network algorithm. In units of 4 bytes. Only used for non-simulated p-values larger than 2×2 tables. This also increases the internal stack size which allows larger problems to be solved, sometimes needing hours. In such cases, <code>simulate.p.values = TRUE</code> may be more reasonable.
<code>alternative</code>	indicates the alternative hypothesis and must be one of "two.sided", "greater" or "less". You can specify just the initial letter. Only used in the 2×2 case.
<code>conf.int</code>	logical indicating if a confidence interval for the odds ratio in a 2×2 table should be computed (and returned).
<code>conf.level</code>	confidence level for the returned confidence interval. Only used in the 2×2 case and if <code>conf.int = TRUE</code> .
<code>simulate.p.value</code>	a logical indicating whether to compute p-values by Monte Carlo simulation, in larger than 2×2 tables.
<code>B</code>	an integer specifying the number of replicates used in the Monte Carlo test when <code>simulate.p.value</code> is true.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>...</code>	Other arguments passed to the function <code>fisher_test()</code> .
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .

Value

return a data frame with some the following columns:

- `group`: the categories in the row-wise proportion tests.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the used statistical test.
- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: an estimate of the odds ratio. Only present in the 2 by 2 case.
- `alternative`: a character string describing the alternative hypothesis.
- `conf.low`, `conf.high`: a confidence interval for the odds ratio. Only present in the 2 by 2 case and if argument `conf.int = TRUE`.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `fisher_test()`: performs Fisher's exact test for testing the null of independence of rows and columns in a contingency table with fixed marginals. Wrapper around the function `fisher.test()`.
- `pairwise_fisher_test()`: pairwise comparisons between proportions, a post-hoc tests following a significant Fisher's exact test of homogeneity for 2xc design.
- `row_wise_fisher_test()`: performs row-wise Fisher's exact test of count data, a post-hoc tests following a significant chi-square test of homogeneity for rx2 contingency table. The test is conducted for each category (row).

See Also

The Datanovia tutorial: Fisher's Exact Test in R.

Examples

```
# Comparing two proportions
#####
# Data: frequencies of smokers between two groups
xtab <- as.table(rbind(c(490, 10), c(400, 100)))
dimnames(xtab) <- list(
  group = c("grp1", "grp2"),
  smoker = c("yes", "no")
)
xtab
# compare the proportion of smokers
fisher_test(xtab, detailed = TRUE)

# Homogeneity of proportions between groups
#####
# H0: the proportion of smokers is similar in the four groups
# Ha: this proportion is different in at least one of the populations.
#
# Data preparation
grp.size <- c( 106, 113, 156, 102 )
smokers <- c( 50, 100, 139, 80 )
no.smokers <- grp.size - smokers
xtab <- as.table(rbind(
  smokers,
  no.smokers
))
dimnames(xtab) <- list(
  Smokers = c("Yes", "No"),
  Groups = c("grp1", "grp2", "grp3", "grp4")
)
xtab

# Compare the proportions of smokers between groups
fisher_test(xtab, detailed = TRUE)

# Pairwise comparison between groups
```

```

pairwise_fisher_test(xtab)

# Pairwise proportion tests
#####
# Data: Titanic
xtab <- as.table(rbind(
  c(122, 167, 528, 673),
  c(203, 118, 178, 212)
))
dimnames(xtab) <- list(
  Survived = c("No", "Yes"),
  Class = c("1st", "2nd", "3rd", "Crew")
)
xtab
# Compare the proportion of survived between groups
pairwise_fisher_test(xtab)

# Row-wise proportion tests
#####
# Data: Titanic
xtab <- as.table(rbind(
  c(180, 145), c(179, 106),
  c(510, 196), c(862, 23)
))
dimnames(xtab) <- list(
  Class = c("1st", "2nd", "3rd", "Crew"),
  Gender = c("Male", "Female")
)
xtab
# Compare the proportion of males and females in each category
row_wise_fisher_test(xtab)

# A r x c table Agresti (2002, p. 57) Job Satisfaction
Job <- matrix(c(1,2,1,0, 3,3,6,1, 10,10,14,9, 6,7,12,11), 4, 4,
  dimnames = list(income = c("< 15k", "15-25k", "25-40k", "> 40k"),
    satisfaction = c("VeryD", "LittleD", "ModerateS", "VeryS")))
fisher_test(Job)
fisher_test(Job, simulate.p.value = TRUE, B = 1e5)

```

fligner_test

Fligner-Killeen Test

Description

Provides a pipe-friendly framework to perform the Fligner-Killeen test, a non-parametric (rank-based) test of the homogeneity of group variances. It is robust against departures from normality and is a useful alternative to `levene_test()`. Wrapper around the function `fligner.test()`.

See the Datanovia tutorial [Homogeneity of Variance Test in R](#) for a worked walkthrough.

Usage

```
fligner_test(data, formula, ...)
```

Arguments

data	a data.frame containing the variables in the formula.
formula	a formula of the form $x \sim \text{group}$ where x is a numeric variable giving the data values and group is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
...	other arguments to be passed to the function <code>fligner.test</code> .

Value

return a data frame with the following columns:

- `.y.`: the y variable used in the test.
- `n`: sample count.
- `statistic`: the Fligner-Killeen test statistic (a chi-squared statistic) used to compute the p-value.
- `df`: the degrees of freedom.
- `p`: p-value.
- `method`: the statistical test used to compare groups.

See Also

[levene_test](#) The Datanovia tutorial: [Homogeneity of Variance Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# Fligner-Killeen test
#::::::::::::::::::::::::::::::::::::
df %>% fligner_test(len ~ dose)

# Grouped data
df %>%
  group_by(supp) %>%
  fligner_test(len ~ dose)
```

freq_table	<i>Compute Frequency Table</i>
------------	--------------------------------

Description

compute frequency table.

Usage

```
freq_table(data, ..., vars = NULL, na.rm = TRUE)
```

Arguments

data	a data frame
...	One or more unquoted expressions (or variable names) separated by commas. Used to specify variables of interest.
vars	optional character vector containing variable names.
na.rm	logical value. If TRUE (default), remove missing values in the variables used to create the frequency table.

Value

a data frame

Examples

```
data("ToothGrowth")
ToothGrowth %>% freq_table(supp, dose)

# Grouped data: frequencies are computed within each group
ToothGrowth %>% group_by(supp) %>% freq_table(dose)
```

friedman_conover_test	<i>Conover's All-Pairs Comparisons Test for Friedman Rank Sums</i>
-----------------------	--

Description

Performs Conover's all-pairs comparison test (also known as the Durbin-Conover test) for a two-way balanced complete block design, following a significant Friedman rank sum test. It is the repeated-measures analogue of Conover's test for the Kruskal-Wallis design: the within-block ranks are compared pairwise using the pooled rank variance and the statistic is referred to a t -distribution with $(b-1)(k-1)$ degrees of freedom (b blocks, k treatments). It should only be used as a post-hoc procedure when the Friedman test is itself significant (Conover, 1999).

If a reference group is specified (via `ref.group`), then each of the remaining treatments is compared only to the reference (control) treatment, and the p-value adjustment for multiple comparisons is computed over only these $k-1$ comparisons (as for [dunn_test\(\)](#)).

See the Datanovia tutorial [Friedman Test in R](#) for a worked walkthrough.

Usage

```

friedman_conover_test(
  data,
  formula,
  p.adjust.method = "holm",
  ref.group = NULL,
  detailed = FALSE
)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form $a \sim b \mid c$, where a (numeric) is the dependent variable name; b is the within-subjects factor variable (the treatment); and c is the column name containing the individuals/subjects (block) identifier. Should be unique per individual.
<code>p.adjust.method</code>	method to adjust p-values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". Default is "holm".
<code>ref.group</code>	a character string specifying the reference treatment. If specified, each of the treatment levels is compared to the reference (control), and the p-value adjustment is computed over only these comparisons.
<code>detailed</code>	logical value. If TRUE, returns the rank-sum estimate and the test method in the output.

Details

For a balanced complete block design with b blocks and k treatments, the observations within each block are ranked. Let R_j be the sum of the within-block ranks for treatment j and let $A = \sum r^2$ be the sum of the squared within-block ranks. The pairwise statistic for treatments i and j is

$$t_{ij} = \frac{R_i - R_j}{\sqrt{\frac{2(bA - \sum_j R_j^2)}{(b-1)(k-1)}}}$$

referred to a t -distribution with $(b-1)(k-1)$ degrees of freedom. This is the Conover (1999) post-hoc, also known as the Durbin-Conover test. In the returned table each row is oriented with $i = \text{group2}$ and $j = \text{group1}$: estimate is $R_{\text{group2}} - R_{\text{group1}}$ and statistic carries its sign, the same convention as `conover_test()` and `dunn_test()`.

The p-values match `PMCMRplus::frdAllPairsConoverTest()`. That function reports the t statistic for the reversed comparison, so its sign is the opposite of the one returned here; the magnitude is the same.

Value

return a data frame with some of the following columns:

- `.y.`: the y (outcome) variable used in the test.
- `group1, group2`: the compared treatments in the pairwise tests.
- `n1, n2`: the number of blocks (subjects) contributing to each treatment.
- `estimate`: the rank-sum difference.
- `estimate1, estimate2`: the rank sums of the two treatments, respectively.
- `statistic`: Test statistic (t-value) used to compute the p-value.
- `df`: degrees of freedom $((b - 1)(k - 1))$.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the statistical test used to compare groups.
- `p.adj.signif`: the significance level of the adjusted p-values.

The **returned object** has an **attribute** called **args**, which is a list holding the test arguments.

References

Conover, W. J. (1999) Practical Nonparametric Statistics, 3rd edition. Wiley.

See Also

[friedman_test](#), [friedman_nemenyi_test](#), [friedman_effsize](#) The Datanovia tutorial: [Friedman Test in R](#).

Examples

```
# A balanced complete block design: 3 treatments measured on 6 subjects
df <- data.frame(
  id      = factor(rep(1:6, 3)),
  treatment = factor(rep(c("A", "B", "C"), each = 6)),
  score    = c(4, 6, 3, 5, 4, 5, 7, 8, 6, 7, 9, 6, 6, 9, 7, 8, 8, 9)
)

# Omnibus Friedman test
df %>% friedman_test(score ~ treatment | id)

# Conover (Durbin-Conover) all-pairs post-hoc
df %>% friedman_conover_test(score ~ treatment | id)

# Comparison against a reference (control) treatment
df %>% friedman_conover_test(score ~ treatment | id, ref.group = "A")
```

friedman_effsize	<i>Friedman Test Effect Size (Kendall's W Value)</i>
------------------	--

Description

Compute the effect size estimate (referred to as w) for Friedman test: $W = X^2/N(K-1)$; where W is the Kendall's W value; X^2 is the Friedman test statistic value; N is the sample size. k is the number of measurements per subject.

The Kendall's W coefficient assumes the value from 0 (indicating no relationship) to 1 (indicating a perfect relationship).

Kendalls uses the Cohen's interpretation guidelines of $0.1 - < 0.3$ (small effect), $0.3 - < 0.5$ (moderate effect) and ≥ 0.5 (large effect)

Confidence intervals are calculated by bootstrap.

See the Datanovia tutorial [Friedman Test in R](#) for a worked walkthrough.

Usage

```
friedman_effsize(
  data,
  formula,
  ci = FALSE,
  conf.level = 0.95,
  ci.type = "perc",
  nboot = 1000,
  ...,
  boot.parallel = getOption("boot.parallel", "no"),
  boot.ncpus = getOption("boot.ncpus", 1L)
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form $a \sim b \mid c$, where a (numeric) is the dependent variable name; b is the within-subjects factor variables; and c (factor) is the column name containing individuals/subjects identifier. Should be unique per individual.
<code>ci</code>	If TRUE, returns confidence intervals by bootstrap. May be slow.
<code>conf.level</code>	The level for the confidence interval.
<code>ci.type</code>	The type of confidence interval to use. Can be any of "norm", "basic", "perc", or "bca". Passed to <code>boot::boot.ci</code> .
<code>nboot</code>	The number of replications to use for bootstrap.
<code>...</code>	other arguments passed to the function <code>friedman.test()</code>

<code>boot.parallel</code>	The type of parallel operation to be used when computing the bootstrap confidence interval. Allowed values are "no" (default), "multicore" and "snow". Passed to <code>boot()</code> . Defaults to <code>getOption("boot.parallel", "no")</code> , so it can also be set globally with <code>options(boot.parallel = "multicore")</code> . Only used when <code>ci = TRUE</code> .
<code>boot.ncpus</code>	Integer. The number of processes to be used in the parallel bootstrap. Defaults to <code>getOption("boot.ncpus", 1L)</code> . Note that <code>boot.parallel</code> has no effect unless <code>boot.ncpus > 1</code> . Only used when <code>ci = TRUE</code> .

Value

return a data frame with some of the following columns:

- `.y.:` the y variable used in the test.
- `n:` Sample counts.
- `effsize:` estimate of the effect size.
- `magnitude:` magnitude of effect size.
- `conf.low, conf.high:` lower and upper bound of the effect size confidence interval.

References

Maciej Tomczak and Ewa Tomczak. The need to report effect size estimates revisited. An overview of some recommended measures of effect size. *Trends in Sport Sciences*. 2014; 1(21):19-25.

See Also

The Datanovia tutorial: [Friedman Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth %>%
  filter(supp == "VC") %>%
  mutate(id = rep(1:10, 3))
head(df)

# Friedman test effect size
#::::::::::::::::::::::::::::::::::::
df %>% friedman_effsize(len ~ dose | id)
```

`friedman_nemenyi_test` *Nemenyi Post-Hoc Test for Friedman Rank Sums*

Description

Performs the Nemenyi (Wilcoxon-Nemenyi-McDonald-Thompson) all-pairs post-hoc test for a two-way balanced complete block design, following a significant Friedman rank sum test. The treatment rank sums are compared pairwise and the test statistic is referred to the studentized range distribution, which already accounts for the multiplicity of the all-pairs comparisons (so, as for `tukey_hsd()`, there is no separate p-value adjustment step). It should only be used as a post-hoc procedure when the Friedman test is itself significant.

The Nemenyi test is the rank-based, repeated-measures analogue of Tukey's HSD. Unlike `friedman_conover_test()` (the Durbin-Conover test), it does not borrow the residual rank variance and is therefore more conservative.

See the Datanovia tutorial [Friedman Test in R](#) for a worked walkthrough.

Usage

```
friedman_nemenyi_test(data, formula, detailed = FALSE)
```

Arguments

<code>data</code>	a <code>data.frame</code> containing the variables in the formula.
<code>formula</code>	a formula of the form <code>a ~ b c</code> , where <code>a</code> (numeric) is the dependent variable name; <code>b</code> is the within-subjects factor variable (the treatment); and <code>c</code> is the column name containing the individuals/subjects (block) identifier. Should be unique per individual.
<code>detailed</code>	logical value. If <code>TRUE</code> , returns the rank-sum estimate and the test method in the output.

Details

For a balanced complete block design with b blocks and k treatments, the observations within each block are ranked. Let R_j be the sum of the within-block ranks for treatment j . The pairwise statistic for treatments i and j is

$$q_{ij} = \frac{|R_i - R_j|}{\sqrt{b k (k + 1) / 12}}$$

and the p-value is obtained from the studentized range distribution with k groups and infinite degrees of freedom.

The p-values match `PMCMRplus::frdAllPairsNemenyiTest()`. That function reports the magnitude of the statistic; the value returned here carries the sign of the rank-sum difference between the two groups, so the two agree only where that difference is positive.

Value

return a data frame with some of the following columns:

- `.y.`: the y (outcome) variable used in the test.
- `group1, group2`: the compared treatments in the pairwise tests.
- `n1, n2`: the number of blocks (subjects) contributing to each treatment.
- `estimate`: the rank-sum difference.
- `estimate1, estimate2`: the rank sums of the two treatments, respectively.
- `statistic`: the studentized-range test statistic.
- `p.adj`: the p-value (already adjusted for multiple comparisons via the studentized range distribution).
- `method`: the statistical test used to compare groups.
- `p.adj.signif`: the significance level of the adjusted p-values.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

References

- Nemenyi, P. (1963) Distribution-free Multiple Comparisons. PhD Thesis, Princeton University.
- Hollander, M., Wolfe, D. A. (1973) Nonparametric Statistical Methods. Wiley.

See Also

[friedman_test](#), [friedman_conover_test](#), [friedman_effsize](#) The Datanovia tutorial: [Friedman Test in R](#).

Examples

```
# A balanced complete block design: 3 treatments measured on 6 subjects
df <- data.frame(
  id      = factor(rep(1:6, 3)),
  treatment = factor(rep(c("A", "B", "C"), each = 6)),
  score   = c(4, 6, 3, 5, 4, 5, 7, 8, 6, 7, 9, 6, 6, 9, 7, 8, 8, 9)
)

# Omnibus Friedman test
df %>% friedman_test(score ~ treatment | id)

# Nemenyi all-pairs post-hoc
df %>% friedman_nemenyi_test(score ~ treatment | id)
```

friedman_test	<i>Friedman Rank Sum Test</i>
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Description

Provides a pipe-friendly framework to perform a Friedman rank sum test, which is the non-parametric alternative to the one-way repeated measures ANOVA test. Wrapper around the function `friedman.test()`.

See the Datanovia tutorial [Friedman Test in R](#) for a worked walkthrough.

Usage

```
friedman_test(data, formula, ...)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>a ~ b c</code> , where <code>a</code> (numeric) is the dependent variable name; <code>b</code> is the within-subjects factor variables; and <code>c</code> (factor) is the column name containing individuals/subjects identifier. Should be unique per individual.
<code>...</code>	other arguments to be passed to the function <code>friedman.test</code> .

Value

return a data frame with the following columns:

- `.y.:` the y (dependent) variable used in the test.
- `n:` sample count.
- `statistic:` the value of Friedman's chi-squared statistic, used to compute the p-value.
- `p:` p-value.
- `method:` the statistical test used to compare groups.

See Also

The Datanovia tutorial: [Friedman Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth %>%
  filter(supp == "VC") %>%
  mutate(id = rep(1:10, 3))
head(df)

# Friedman rank sum test
#::::::::::::::::::::::::::::::::::::
```

```
df %>% friedman_test(len ~ dose | id)
```

games_howell_test	<i>Games Howell Post-hoc Tests</i>
-------------------	------------------------------------

Description

Performs Games-Howell test, which is used to compare all possible combinations of group differences when the assumption of homogeneity of variances is violated. This post hoc test provides confidence intervals for the differences between group means and shows whether the differences are statistically significant.

The test is based on Welch's degrees of freedom correction and uses Tukey's studentized range distribution for computing the p-values. The test compares the difference between each pair of means with appropriate adjustment for the multiple testing. So there is no need to apply additional p-value corrections.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
games_howell_test(
  data,
  formula,
  conf.level = 0.95,
  detailed = FALSE,
  effect.size = FALSE
)
```

Arguments

data	a data.frame containing the variables in the formula.
formula	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
conf.level	confidence level of the interval.
detailed	logical value. Default is FALSE. If TRUE, a detailed result is shown.
effect.size	logical. Default is FALSE. If TRUE, a <code>cohens.d</code> column and its magnitude are added. Because Games-Howell assumes unequal variances, this is the Welch (averaged-variance) Cohen's d, $\text{estimate} / \sqrt{(s_1^2 + s_2^2) / 2}$, oriented like the reported mean difference so the two never disagree in sign.

Details

The Games-Howell method is an improved version of the Tukey-Kramer method and is applicable in cases where the equivalence of variance assumption is violated. It is a t-test using Welch's degree of freedom. This method uses a strategy for controlling the type I error for the entire comparison and is known to maintain the preset significance level even when the size of the sample is different. However, the smaller the number of samples in each group, the it is more tolerant the type I error control. Thus, this method can be applied when the number of samples is six or more.

Because the test relies on Welch's variance correction, comparisons involving a group with zero variance (constant values) or undefined variance (a single observation) can be undefined; such pairs are returned as NA (with a warning) while all other comparisons are computed as usual.

Value

return a data frame with some of the following columns:

- `.y.`: the y (outcome) variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n1, n2`: Sample counts.
- `estimate, conf.low, conf.high`: mean difference and its confidence intervals.
- `statistic`: Test statistic (t-value) used to compute the p-value.
- `df`: degrees of freedom calculated using Welch's correction.
- `p.adj`: adjusted p-value using Tukey's method.
- `method`: the statistical test used to compare groups.
- `p.adj.signif`: the significance level of p-values.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

References

- Aaron Schlege, <https://rpubs.com/aaronsc32/games-howell-test>.
- Sangseok Lee, Dong Kyu Lee. What is the proper way to apply the multiple comparison test?. Korean J Anesthesiol. 2018;71(5):353-360.

See Also

The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Simple test
ToothGrowth %>% games_howell_test(len ~ dose)

# Grouped data
ToothGrowth %>%
  group_by(supp) %>%
  games_howell_test(len ~ dose)
```

get_comparisons	<i>Create a List of Possible Comparisons Between Groups</i>
-----------------	---

Description

Create a list of possible pairwise comparisons between groups. If a reference group is specified, only comparisons against reference will be kept.

Usage

```
get_comparisons(data, variable, ref.group = NULL)
```

Arguments

data	a data frame
variable	the grouping variable name. Can be unquoted.
ref.group	a character string specifying the reference group. Can be unquoted. If numeric, then it should be quoted. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If ref.group = "all", pairwise comparisons are performed between each grouping variable levels against all (i.e. basemean).

Value

a list of all possible pairwise comparisons.

Examples

```
# All possible pairwise comparisons
ToothGrowth %>%
  get_comparisons("dose")

# Comparisons against reference groups
ToothGrowth %>%
  get_comparisons("dose", ref.group = "0.5")

# Comparisons against all (basemean)
ToothGrowth %>%
  get_comparisons("dose", ref.group = "all")
```

`get_mode`*Compute Mode*

Description

Compute the mode in a given vector. Mode is the most frequent value.

Usage

```
get_mode(x)
```

Arguments

`x` a vector. Can be numeric, factor or character vector.

Examples

```
# Mode of numeric vector
x <- c(1:5, 6, 6, 7:10)
get_mode(x)

# Bimodal
x <- c(1:5, 6, 6, 7, 8, 9, 9, 10)
get_mode(x)

# No mode
x <- c(1, 2, 3, 4, 5)
get_mode(x)

# Nominal vector
fruits <- c(rep("orange", 10), rep("apple", 5), rep("lemon", 2))
get_mode(fruits)
```

`get_pwc_label`*Extract Label Information from Statistical Tests*

Description

Extracts label information from statistical tests. Useful for labelling plots with test outputs.

See the Datanovia tutorial [P-values from Tests on ggplots in R \(rstatix\)](#) for a worked walkthrough.

Usage

```
get_pwc_label(stat.test, type = c("expression", "text"))
```

```
get_test_label(
  stat.test,
  description = NULL,
  p.col = "p",
  type = c("expression", "text"),
  correction = c("auto", "GG", "HF", "none"),
  row = NULL,
  detailed = FALSE,
  style = c("classic", "apa")
)
```

```
create_test_label(
  statistic.text,
  statistic,
  p,
  parameter = NA,
  description = NULL,
  n = NA,
  effect.size = NA,
  effect.size.text = NA,
  type = c("expression", "text"),
  detailed = FALSE,
  style = c("classic", "apa"),
  effect.size.ci = NA,
  effect.size.bounded = TRUE,
  effect.size.ci.level = 0.95
)
```

```
get_n(stat.test)
```

```
get_description(stat.test)
```

Arguments

<code>stat.test</code>	statistical test results returned by <code>rstatix</code> functions.
<code>type</code>	the label type. Can be one of "text" and "expression". Partial match allowed. If you want to add the label onto a <code>ggplot</code> , it might be useful to specify <code>type = "expresion"</code> .
<code>description</code>	the test description used as the prefix of the label. Examples of description are "ANOVA", "Two Way ANOVA". To remove the default description, specify <code>description = NULL</code> . If missing, we'll try to guess the statistical test default description.
<code>p.col</code>	character specifying the column containing the p-value. Default is "p", can be "p.adj".

correction	character, considered only in the case of ANOVA test. Which sphericity correction of the degrees of freedom should be reported for the within-subject factors (repeated measures). The default is set to "GG" corresponding to the Greenhouse-Geisser correction. Possible values are "GG", "HF" (i.e., Hyunh-Feldt correction), "none" (i.e., no correction) and "auto" (apply automatically GG correction if the sphericity assumption is not for within-subject design).
row	numeric, the row index to be considered. If NULL, the last row is automatically considered for ANOVA test.
detailed	logical value. If TRUE, returns detailed label.
style	the label style. Either "classic" (default, the historical format) or "apa" for an APA-7 in-text statistical report: the leading test-name label is dropped, the p-value is formatted APA-style ($p = .023$, $p < .001$), and the effect size is shown with its confidence interval when the object carries one (e.g. <code>anova_test(ci =)</code>), otherwise as a point estimate, otherwise omitted. Bounded effect sizes (η^2 , r , Cliff's δ , rank-biserial) drop the leading zero; Cohen's d keeps it.
statistic.text	character specifying the test statistic. For example <code>statistic.text = "F"</code> (for ANOVA test); <code>statistic.text = "t"</code> (for t-test).
statistic	the numeric value of a statistic.
p	the p-value of the test.
parameter	string containing the degree of freedom (if exists). Default is NA to accommodate non-parametric tests. For example <code>parameter = "1,9"</code> (for ANOVA test. Two parameters exist: DFn and DFd); <code>sparameter = "9"</code> (for t-test).
n	sample count, example: $n = 10$.
effect.size	the effect size value
effect.size.text	a character specifying the relevant effect size. For example, for Cohens d statistic, <code>effect.size.text = "d"</code> . You can also use plotmath expression as follow <code>quote(italic("d"))</code> .
effect.size.ci	a length-two numeric <code>c(low, high)</code> giving the effect size's confidence interval, appended after the effect size when <code>style = "apa"</code> . Defaults to NA (no interval).
effect.size.bounded	logical. Whether the effect size lies in $[-1, 1]$ (e.g. η^2 , r , Cliff's δ); used by <code>style = "apa"</code> to decide whether to drop the leading zero. Defaults to TRUE; set FALSE for Cohen's d .
effect.size.ci.level	the confidence level <code>effect.size.ci</code> was computed at, shown before the interval (e.g. 95% CI). Defaults to 0.95.

Value

a text label or an expression to pass to a plotting function.

Functions

- `get_pwc_label()`: Extract label from pairwise comparisons.

- `get_test_label()`: Extract labels for statistical tests.
- `create_test_label()`: Create labels from user specified test results.
- `get_n()`: Extracts sample counts (n) from an rstatix test outputs. Returns a numeric vector.
- `get_description()`: Extracts the description of an rstatix test outputs. Returns a character vector.

References

American Psychological Association (2020). *Publication Manual of the American Psychological Association* (7th ed.).

See Also

The Datanovia tutorial: [P-values from Tests on ggplots in R \(rstatix\)](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# One-way ANOVA test
#::::::::::::::::::::::::::::::::::::
anov <- df %>% anova_test(len ~ dose)
get_test_label(anov, detailed = TRUE, type = "text")

# Two-way ANOVA test
#::::::::::::::::::::::::::::::::::::
anov <- df %>% anova_test(len ~ supp*dose)
get_test_label(anov, detailed = TRUE, type = "text",
  description = "Two Way ANOVA")

# Kruskal-Wallis test
#::::::::::::::::::::::::::::::::::::
kruskal<- df %>% kruskal_test(len ~ dose)
get_test_label(kruskal, detailed = TRUE, type = "text")

# Wilcoxon test
#::::::::::::::::::::::::::::::::::::
# Unpaired test
wilcox <- df %>% wilcox_test(len ~ supp)
get_test_label(wilcox, detailed = TRUE, type = "text")
# Paired test
wilcox <- df %>% wilcox_test(len ~ supp, paired = TRUE)
get_test_label(wilcox, detailed = TRUE, type = "text")

# T test
#::::::::::::::::::::::::::::::::::::
ttest <- df %>% t_test(len ~ dose)
```

```

get_test_label(ttest, detailed = TRUE, type = "text")

# Pairwise comparisons labels
#::::::::::::::::::::::::::::::::::::
get_pwc_label(ttest, type = "text")

# Create test labels
#::::::::::::::::::::::::::::::::::::
create_test_label(
  statistic.text = "F", statistic = 71.82,
  parameter = "4, 294",
  p = "<0.0001",
  description = "ANOVA",
  type = "text"
)

# Extract infos
#::::::::::::::::::::::::::::::::::::
stat.test <- df %>% t_test(len ~ dose)
get_n(stat.test)
get_description(stat.test)

```

get_summary_stats	<i>Compute Summary Statistics</i>
-------------------	-----------------------------------

Description

Compute summary statistics for one or multiple numeric variables.

See the Datanovia tutorial [Descriptive Statistics in R](#) for a worked walkthrough.

Usage

```

get_summary_stats(
  data,
  ...,
  type = c("full", "common", "robust", "five_number", "mean_sd", "mean_se", "mean_ci",
    "median_iqr", "median_mad", "quantile", "mean", "median", "min", "max"),
  show = NULL,
  probs = seq(0, 1, 0.25),
  digits = 3
)

```

Arguments

<code>data</code>	a data frame
<code>...</code>	(optional) One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest. If no variable is specified, then the summary statistics of all numeric variables in the data frame is computed.
<code>type</code>	type of summary statistics. Possible values include: "full", "common", "robust", "five_number", "mean_sd", "mean_se", "mean_ci", "median_iqr", "median_mad", "quantile", "mean", "median", "min", "max"
<code>show</code>	a character vector specifying the summary statistics you want to show. Example: <code>show = c("n", "mean", "sd")</code> . This is used to filter the output after computation. It can additionally include "skewness" and/or "kurtosis" (e.g. <code>show = c("mean", "sd", "skewness", "kurtosis")</code>); these two are computed on demand and are not part of any default type.
<code>probs</code>	numeric vector of probabilities with values in [0,1]. Used only when <code>type = "quantile"</code> .
<code>digits</code>	integer indicating the number of decimal places to round the summary statistics to. Default is 3. Increase it when summarizing very small values that would otherwise round to 0.

Value

A data frame containing descriptive statistics, such as:

- **n**: the number of individuals
- **min**: minimum
- **max**: maximum
- **median**: median
- **mean**: mean
- **q1, q3**: the first and the third quartile, respectively.
- **iqr**: interquartile range
- **mad**: median absolute deviation (see ?MAD)
- **sd**: standard deviation of the mean
- **se**: standard error of the mean
- **ci**: 95 percent confidence interval of the mean

When requested through `show`, the output can also contain:

- **skewness**: bias-corrected sample skewness
- **kurtosis**: bias-corrected sample excess kurtosis (0 for a normal distribution).

Both use the type-2 (bias-corrected) estimator, matching e1071 with `type = 2`: $\text{skewness} = g_1 \sqrt{n(n-1)} / (n-2)$ and $\text{kurtosis} = [(n+1)g_2 + 6](n-1) / [(n-2)(n-3)]$, where $g_1 = m_3 / m_2^{1.5}$ and $g_2 = m_4 / m_2^2 - 3$. Skewness is NA for $n < 3$ and kurtosis for $n < 4$.

See Also

[rstatix-programming](#) for selecting columns by names held in strings (!!, {{ }}, vars=, all_of()).
 The Datanovia tutorial: [Descriptive Statistics in R](#).

Examples

```
# Full summary statistics
data("ToothGrowth")
ToothGrowth %>% get_summary_stats(len)

# Summary statistics of grouped data
# Show only common summary
ToothGrowth %>%
  group_by(dose, supp) %>%
  get_summary_stats(len, type = "common")

# Robust summary statistics
ToothGrowth %>% get_summary_stats(len, type = "robust")

# Five number summary statistics
ToothGrowth %>% get_summary_stats(len, type = "five_number")

# Compute only mean and sd
ToothGrowth %>% get_summary_stats(len, type = "mean_sd")

# Compute full summary statistics but show only mean, sd, median, iqr
ToothGrowth %>%
  get_summary_stats(len, show = c("mean", "sd", "median", "iqr"))

# Include skewness and kurtosis (computed on demand via show)
ToothGrowth %>%
  get_summary_stats(len, show = c("mean", "sd", "skewness", "kurtosis"))
```

get_y_position

Autocompute P-value Positions For Plotting Significance

Description

Compute p-value x and y positions for plotting significance levels.

See the Datanovia tutorials [P-values from Tests on ggplots in R \(rstatix\)](#) and [Auto P-values in ggplot with geom_pwc \(ggpubr\)](#) for worked walkthroughs.

Usage

```
get_y_position(
  data,
  formula,
```

```

    fun = "max",
    ref.group = NULL,
    comparisons = NULL,
    step.increase = 0.12,
    y.trans = NULL,
    stack = FALSE,
    scales = c("fixed", "free", "free_y")
)

add_y_position(
  test,
  fun = "max",
  step.increase = 0.12,
  data = NULL,
  formula = NULL,
  ref.group = NULL,
  comparisons = NULL,
  y.trans = NULL,
  stack = FALSE,
  scales = c("fixed", "free", "free_y")
)

add_x_position(
  test,
  x = NULL,
  group = NULL,
  dodge = 0.8,
  scales = c("fixed", "free", "free_y")
)

add_xy_position(
  test,
  x = NULL,
  group = NULL,
  dodge = 0.8,
  stack = FALSE,
  fun = "max",
  step.increase = 0.12,
  scales = c("fixed", "free", "free_y"),
  ...
)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .

fun	summary statistics functions used to compute automatically suitable y positions of p-value labels and brackets. Possible values include: "max", "mean", "mean_sd", "mean_se", "mean_ci", "median", "median_iqr", "median_mad". For example, if fun = "max", the y positions are guessed as follow: • 1. Compute the maximum of each group (groups.maximum) • 2. Use the highest groups maximum as the first bracket y position • 3. Add successively a step increase for remaining bracket y positions. When the main plot is a boxplot, you need the option fun = "max", to have the p-value bracket displayed at the maximum point of the group. In some situations the main plot is a line plot or a barplot showing the mean+/-error bars of the groups, where error can be SE (standard error), SD (standard deviation) or CI (confidence interval). In this case, to correctly compute the bracket y position you need the option fun = "mean_se", etc.
ref.group	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group).
comparisons	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: comparisons = list(c("A", "B"), c("B", "C"))
step.increase	numeric vector with the increase in fraction of total height for every additional comparison to minimize overlap.
y.trans	a function for transforming y axis scale. Value can be log2, log10 and sqrt. Can be also any custom function that can take a numeric vector as input and returns a numeric vector, example: y.trans = function(x){log2(x+1)}
stack	logical. If TRUE, computes y position for a stacked plot. Useful when dealing with stacked bar plots.
scales	Should scales be fixed ("fixed", the default), free ("free"), or free in one dimension ("free_y")?. For the y position, "free" or "free_y" compute the step increase per plot panel (otherwise a global step increase is used). For the x position, when scales = "free" the x positions (xmin/xmax) are computed per facet , so that brackets align correctly when a faceted plot (facet_*(scales = "free")) shows a different set of x-axis levels in each panel ("free_y" keeps the global x positions).
test	an object of class rstatix_test as returned by t_test() , wilcox_test() , sign_test() , tukey_hsd() , dunn_test() .
x	variable on x axis.
group	group variable (legend variable).
dodge	dodge width for grouped ggplot/test. Default is 0.8. Used only when x specified.
...	other arguments to be passed to the function t.test .

Functions

- [get_y_position\(\)](#): compute the p-value y positions
- [add_y_position\(\)](#): add p-value y positions to an object of class rstatix_test
- [add_x_position\(\)](#): compute and add p-value x positions.
- [add_xy_position\(\)](#): compute and add both x and y positions.

Added columns

`add_y_position()` adds `y.position`; `add_x_position()` adds `x`, `xmin`, `xmax`; and `add_xy_position()` adds all of these. A column named `groups` (the pair of compared groups, used internally to position dodged grouped comparisons) is also added. **groups is a reserved internal column**: do not rely on it and avoid using `groups` as one of your own column names in the test object, since it is overwritten here. (It may be renamed to a dotted, less collision-prone name in a future major version.)

See Also

The Datanovia tutorials: [P-values from Tests on ggplots in R \(rstatix\)](#), [Auto P-values in ggplot with geom_pwc \(ggpubr\)](#).

Examples

```
# Data preparation
#::::::::::::::::::::::::::::::::::::
df <- ToothGrowth
df$dose <- as.factor(df$dose)
df$group <- factor(rep(c(1, 2), 30))
head(df)

# Stat tests
#::::::::::::::::::::::::::::::::::::
stat.test <- df %>%
  t_test(len ~ dose)
stat.test

# Add the test into box plots
#::::::::::::::::::::::::::::::::::::
stat.test <- stat.test %>%
  add_y_position()

if(require("ggpubr")){
  ggboxplot(df, x = "dose", y = "len") +
    stat_pvalue_manual(stat.test, label = "p.adj.signif", tip.length = 0.01)
}
```

Description

Detect outliers using boxplot methods. Boxplots are a popular and an easy method for identifying outliers. There are two categories of outlier: (1) outliers and (2) extreme points.

Values above $Q3 + 1.5 \times IQR$ or below $Q1 - 1.5 \times IQR$ are considered as outliers. Values above $Q3 + 3 \times IQR$ or below $Q1 - 3 \times IQR$ are considered as extreme points (or extreme outliers).

Q1 and Q3 are the first and third quartile, respectively. IQR is the interquartile range ($IQR = Q3 - Q1$).

Generally speaking, data points that are labelled outliers in boxplots are not considered as troublesome as those considered extreme points and might even be ignored. Note that, any NA and NaN are automatically removed before the quantiles are computed.

See the Datanovia tutorial [Descriptive Statistics in R](#) for a worked walkthrough.

Usage

```
identify_outliers(data, ..., variable = NULL)
```

```
is_outlier(x, coef = 1.5)
```

```
is_extreme(x)
```

Arguments

data	a data frame
...	One unquoted expressions (or variable name). Used to select a variable of interest. Alternative to the argument <code>variable</code> .
variable	variable name for detecting outliers
x	a numeric vector
coef	coefficient specifying how far the outlier should be from the edge of their box. Possible values are 1.5 (for outlier) and 3 (for extreme points only). Default is 1.5

Value

- `identify_outliers()`. Returns the input data frame with two additional columns: "is.outlier" and "is.extreme", which hold logical values.
- `is_outlier()` and `is_extreme()`. Returns logical vectors.

Functions

- `identify_outliers()`: takes a data frame and extract rows suspected as outliers according to a numeric column. The following columns are added "is.outlier" and "is.extreme".
- `is_outlier()`: detect outliers in a numeric vector. Returns logical vector.
- `is_extreme()`: detect extreme points in a numeric vector. An alias of `is_outlier()`, where `coef = 3`. Returns logical vector.

See Also

The Datanovia tutorial: [Descriptive Statistics in R](#).

Examples

```
# Generate a demo data
set.seed(123)
demo.data <- data.frame(
  sample = 1:20,
  score = c(rnorm(19, mean = 5, sd = 2), 50),
  gender = rep(c("Male", "Female"), each = 10)
)

# Identify outliers according to the variable score
demo.data %>%
  identify_outliers(score)

# Identify outliers by groups
demo.data %>%
  group_by(gender) %>%
  identify_outliers("score")
```

kruskal_effsize

Kruskal-Wallis Effect Size

Description

Compute the effect size for Kruskal-Wallis test as the eta squared based on the H-statistic: $\eta^2[H] = (H - k + 1) / (n - k)$; where H is the value obtained in the Kruskal-Wallis test; k is the number of groups; n is the total number of observations.

The eta-squared estimate assumes values from 0 to 1 and multiplied by 100 indicates the percentage of variance in the dependent variable explained by the independent variable. The interpretation values commonly in published literature are: $0.01 - < 0.06$ (small effect), $0.06 - < 0.14$ (moderate effect) and ≥ 0.14 (large effect).

Note that $\eta^2[H]$ is a bias-corrected estimator, so the raw formula can return a small negative value for a near-null effect (very small H). In that case the estimate is floored to 0, keeping the reported effect size within its valid $[0, 1]$ range.

Confidence intervals are calculated by bootstrap.

See the Datanovia tutorial [Kruskal-Wallis Test in R](#) for a worked walkthrough.

Usage

```
kruskal_effsize(
  data,
  formula,
  ci = FALSE,
  conf.level = 0.95,
  ci.type = "perc",
  nboot = 1000,
  boot.parallel = getOption("boot.parallel", "no"),
```

```

boot.ncpus = getOption("boot.ncpus", 1L),
method = c("eta2", "epsilon2")
)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>ci</code>	If TRUE, returns confidence intervals by bootstrap. May be slow.
<code>conf.level</code>	The level for the confidence interval.
<code>ci.type</code>	The type of confidence interval to use. Can be any of "norm", "basic", "perc", or "bca". Passed to <code>boot::boot.ci</code> .
<code>nboot</code>	The number of replications to use for bootstrap.
<code>boot.parallel</code>	The type of parallel operation to be used when computing the bootstrap confidence interval. Allowed values are "no" (default), "multicore" and "snow". Passed to <code>boot()</code> . Defaults to <code>getOption("boot.parallel", "no")</code> , so it can also be set globally with <code>options(boot.parallel = "multicore")</code> . Only used when <code>ci = TRUE</code> .
<code>boot.ncpus</code>	Integer. The number of processes to be used in the parallel bootstrap. Defaults to <code>getOption("boot.ncpus", 1L)</code> . Note that <code>boot.parallel</code> has no effect unless <code>boot.ncpus > 1</code> . Only used when <code>ci = TRUE</code> .
<code>method</code>	the effect-size metric. Either "eta2" (default) for the bias-corrected eta-squared $\eta^2[H] = (H - k + 1) / (N - k)$, or "epsilon2" for the rank epsilon-squared $H / (N - 1)$ (Tomczak & Tomczak, 2014), which equals <code>effectsize::rank_epsilon_squared()</code> and is not bias-corrected (so a little larger than $\eta^2[H]$).

Value

return a data frame with some of the following columns:

- `.y`: the y variable used in the test.
- `n`: Sample counts.
- `effsize`: estimate of the effect size.
- `magnitude`: magnitude of effect size.
- `conf.low`, `conf.high`: lower and upper bound of the effect size confidence interval.

References

Maciej Tomczak and Ewa Tomczak. The need to report effect size estimates revisited. An overview of some recommended measures of effect size. *Trends in Sport Sciences*. 2014; 1(21):19-25.

<http://imaging.mrc-cbu.cam.ac.uk/statswiki/FAQ/effectSize>

<http://www.psy.gla.ac.uk/~steve/best/effect.html>

See Also

The Datanovia tutorial: [Kruskal-Wallis Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# Kruskal-wallis rank sum test
#::::::::::::::::::::::::::::::::::::
df %>% kruskal_effsize(len ~ dose)

# Grouped data
df %>%
  group_by(supp) %>%
  kruskal_effsize(len ~ dose)
```

kruskal_test	<i>Kruskal-Wallis Test</i>
--------------	----------------------------

Description

Provides a pipe-friendly framework to perform Kruskal-Wallis rank sum test. Wrapper around the function `kruskal.test()`.

See the Datanovia tutorial [Kruskal-Wallis Test in R](#) for a worked walkthrough.

Usage

```
kruskal_test(data, formula, ...)
```

Arguments

data	a data.frame containing the variables in the formula.
formula	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
...	other arguments to be passed to the function <code>kruskal.test</code> .

Value

return a data frame with the following columns:

- `.y.`: the y variable used in the test.
- `n`: sample count.
- `statistic`: the kruskal-wallis rank sum statistic used to compute the p-value.
- `p`: p-value.
- `method`: the statistical test used to compare groups.

See Also

The Datanovia tutorial: [Kruskal-Wallis Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# Kruskal-wallis rank sum test
#::::::::::::::::::::::::::::::::::::
df %>% kruskal_test(len ~ dose)

# Grouped data
df %>%
  group_by(supp) %>%
  kruskal_test(len ~ dose)
```

ks_test

Two-Sample Kolmogorov-Smirnov Test

Description

Provides a pipe-friendly framework to perform the two-sample Kolmogorov-Smirnov test, comparing the (empirical) distributions of a numeric variable between two groups. Wrapper around the R base function `ks.test()`.

When the grouping factor contains more than two levels, pairwise Kolmogorov-Smirnov tests are automatically performed, with p-value adjustment.

See the Datanovia tutorial [Normality Test in R](#) for a worked walkthrough.

Usage

```
ks_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  alternative = "two.sided",
  exact = NULL,
  detailed = FALSE
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with two or more levels giving the corresponding groups.
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code> .
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable level against all (i.e. basemean).
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>alternative</code>	indicates the alternative hypothesis and must be one of "two.sided" (default), "less", or "greater". You can specify just the initial letter of the value, but the argument name must be given in full. See 'Details' for the meanings of the possible values.
<code>exact</code>	NULL or a logical indicating whether an exact p-value should be computed. See 'Details' for the meaning of NULL.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.

Value

return a data frame with some of the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n1, n2`: sample counts.
- `statistic`: the value of the test statistic D (the maximum difference between the two empirical cumulative distribution functions).
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the statistical test used to compare groups.
- `p.signif, p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `alternative`: the alternative hypothesis.

The **returned object has an attribute called args**, which is a list holding the test arguments.

See Also

`wilcox_test()`, `t_test()` The Datanovia tutorial: [Normality Test in R](#).

Examples

```
# Two-samples test
#::::::::::::::::::::::::::::::::::::
ToothGrowth %>% ks_test(len ~ supp)

# Pairwise comparisons (more than two groups)
#::::::::::::::::::::::::::::::::::::
ToothGrowth %>% ks_test(len ~ dose)

# Comparison against a reference group
#::::::::::::::::::::::::::::::::::::
ToothGrowth %>% ks_test(len ~ dose, ref.group = "0.5")
```

levene_test

Levene's Test

Description

Provide a pipe-friendly framework to easily compute Levene's test for homogeneity of variance across groups.

Wrapper around the function `leveneTest()`, which can additionally handles a grouped data.

See the Datanovia tutorial [Homogeneity of Variance Test in R](#) for a worked walkthrough.

Usage

```
levene_test(data, formula, center = median)
```

Arguments

data	a data frame for evaluating the formula or a model
formula	a formula
center	The name of a function to compute the center of each group; mean gives the original Levene's test; the default, median, provides a more robust test.

Value

a data frame with the following columns: df1, df2 (df.residual), statistic and p.

See Also

The Datanovia tutorial: [Homogeneity of Variance Test in R](#).

Examples

```
# Prepare the data
data("ToothGrowth")
df <- ToothGrowth
df$dose <- as.factor(df$dose)
# Compute Levene's Test
df %>% levene_test(len ~ dose)

# Grouped data
df %>%
  group_by(supp) %>%
  levene_test(len ~ dose)
```

`mahalanobis_distance` *Compute Mahalanobis Distance and Flag Multivariate Outliers*

Description

Pipe-friendly wrapper around to the function `mahalanobis()`, which returns the squared Mahalanobis distance of all rows in `x`. Compared to the base function, it automatically flags multivariate outliers.

Mahalanobis distance is a common metric used to identify multivariate outliers. The larger the value of Mahalanobis distance, the more unusual the data point (i.e., the more likely it is to be a multivariate outlier).

The distance tells us how far an observation is from the center of the cloud, taking into account the shape (covariance) of the cloud as well.

To detect outliers, the calculated Mahalanobis distance is compared against a chi-square (X^2) distribution with degrees of freedom equal to the number of dependent (outcome) variables and an alpha level of 0.001.

The threshold to declare a multivariate outlier is determined using the function `qchisq(0.999, df)`, where `df` is the degree of freedom (i.e., the number of dependent variable used in the computation).

See the Datanovia tutorial [Normality Test in R](#) for a worked walkthrough.

Usage

```
mahalanobis_distance(data, ...)
```

Arguments

<code>data</code>	a data frame. Columns are variables.
<code>...</code>	One unquoted expressions (or variable name). Used to select a variable of interest. Can be also used to ignore a variable that are not needed for the computation. For example specify <code>-id</code> to ignore the <code>id</code> column.

Value

Returns the input data frame with two additional columns: 1) "mahal.dist": Mahalanobis distance values; and 2) "is.outlier": logical values specifying whether a given observation is a multivariate outlier

See Also

The Datanovia tutorial: [Normality Test in R](#).

Examples

```
# Compute mahalanobis distance and flag outliers if any
iris %>%
  doo(~mahalanobis_distance(.))

# Compute distance by groups and filter outliers
iris %>%
  group_by(Species) %>%
  doo(~mahalanobis_distance(.)) %>%
  filter(is.outlier == TRUE)
```

make_clean_names	<i>Make Clean Names</i>
------------------	-------------------------

Description

Pipe-friendly function to make syntactically valid names out of character vectors.

Usage

```
make_clean_names(data)
```

Arguments

data a data frame or vector

Value

a data frame or a vector depending on the input data

Examples

```
# Vector
make_clean_names(c("a and b", "a-and-b"))
make_clean_names(1:10)

# data frame
df <- data.frame(
```

```
`a and b` = 1:4,
`c and d` = 5:8,
  check.names = FALSE
)
df
make_clean_names(df)
```

mcnemar_test

*McNemar's Chi-squared Test for Count Data***Description**

Performs McNemar chi-squared test to compare paired proportions.

Wrappers around the R base function `mcnemar.test()`, but provide pairwise comparisons between multiple groups

See the Datanovia tutorial [McNemar's Test in R](#) for a worked walkthrough.

Usage

```
mcnemar_test(x, y = NULL, correct = TRUE)
```

```
pairwise_mcnemar_test(
  data,
  formula,
  type = c("mcnemar", "exact"),
  correct = TRUE,
  p.adjust.method = "bonferroni"
)
```

Arguments

<code>x</code>	either a two-dimensional contingency table in matrix form, or a factor object.
<code>y</code>	a factor object; ignored if <code>x</code> is a matrix.
<code>correct</code>	a logical indicating whether to apply continuity correction when computing the test statistic.
<code>data</code>	a data frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>a ~ b c</code> , where <code>a</code> is the outcome variable name; <code>b</code> is the within-subjects factor variables; and <code>c</code> (factor) is the column name containing individuals/subjects identifier. Should be unique per individual.
<code>type</code>	type of statistical tests used for pairwise comparisons. Allowed values are one of <code>c("mcnemar", "exact")</code> .
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .

Value

return a data frame with the following columns:

- `n`: the number of participants.
- `statistic`: the value of McNemar's statistic.
- `df`: the degrees of freedom of the approximate chi-squared distribution of the test statistic.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the used statistical test.
- `p.signif`: the significance level of p-values.

For `pairwise_mcnemar_test()`, the `statistic` and `df` columns are returned for `type = "mcnemar"` (the default); they are omitted for `type = "exact"`, which is based on an exact binomial test and has no chi-squared statistic or degrees of freedom.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `mcnemar_test()`: performs McNemar's chi-squared test for comparing two paired proportions
- `pairwise_mcnemar_test()`: performs pairwise McNemar's chi-squared test between multiple groups. Could be used for post-hoc tests following a significant Cochran's Q test.

See Also

The Datanovia tutorial: [McNemar's Test in R](#).

Examples

```
# Comparing two paired proportions
# %%%%%%%%%%%
# Data: frequencies of smokers before and after interventions
xtab <- as.table(
  rbind(c(25, 6), c(21,10))
)
dimnames(xtab) <- list(
  before = c("non.smoker", "smoker"),
  after = c("non.smoker", "smoker")
)
xtab

# Compare the proportion of smokers
mcnemar_test(xtab)

# Comparing multiple related proportions
# %%%%%%%%%%%
# Generate a demo data
mydata <- data.frame(
```

```
outcome = c(0,1,1,0,0,1,0,1,1,1,1,1,0,0,1,1,0,1,0,1,1,0,0,1,0,1,1,0,0,1),
treatment = gl(3,1,30,labels=LETTERS[1:3]),
participant = gl(10,3,labels=letters[1:10])
)
mydata$outcome <- factor(
  mydata$outcome, levels = c(1, 0),
  labels = c("success", "failure")
)
# Cross-tabulation
xtabs(~outcome + treatment, mydata)

# Compare the proportion of success between treatments
cochran_qtest(mydata, outcome ~ treatment|participant)

# pairwise comparisons between groups
pairwise_mcnemar_test(mydata, outcome ~ treatment|participant)
```

multinom_test	<i>Exact Multinomial Test</i>
---------------	-------------------------------

Description

Performs an exact multinomial test. Alternative to the chi-square test of goodness-of-fit-test when the sample size is small.

Usage

```
multinom_test(x, p = rep(1/length(x), length(x)), detailed = FALSE)
```

Arguments

- x numeric vector containing the counts.
- p a vector of probabilities of success. The length of p must be the same as the number of groups specified by x, and its elements must be greater than 0 and less than 1.
- detailed logical value. Default is FALSE. If TRUE, a detailed result is shown.

Value

return a data frame containing the p-value and its significance.
The **returned object has an attribute called args**, which is a list holding the test arguments.

See Also

[binom_test](#)

Examples

```
# Data
tulip <- c(red = 81, yellow = 50, white = 27)

# Question 1: are the color equally common ?
#####
# this is a test of homogeneity
res <- multinom_test(tulip)
res

attr(res, "descriptives")

# Pairwise comparisons between groups
pairwise_binom_test(tulip, p.adjust.method = "bonferroni")

# Question 2: comparing observed to expected proportions
#####
# this is a goodness-of-fit test
expected.p <- c(red = 0.5, yellow = 0.33, white = 0.17)
res <- multinom_test(tulip, expected.p)
res
attr(res, "descriptives")

# Pairwise comparisons against a given probabilities
pairwise_binom_test_against_p(tulip, expected.p)
```

omega_squared	<i>Omega Squared for ANOVA</i>
---------------	--------------------------------

Description

Compute (classic, full) omega-squared and partial omega-squared for all terms in a between-subjects ANOVA model. Omega squared is a less-biased alternative to eta squared, estimating the population effect size rather than the sample one.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
omega_squared(model)

partial_omega_squared(model)
```

Arguments

model	an object of class aov or anova (a between-subjects design). Repeated-measures and mixed models are not supported, as for eta_squared() .
-------	---

Value

a named numeric vector of effect sizes, one per model term. A negative point estimate (which can arise for a term with $F < 1$) is floored at 0, since omega squared estimates a non-negative proportion of variance.

Functions

- `omega_squared()`: compute the classic (full) omega squared.
- `partial_omega_squared()`: compute partial omega squared.

References

Olejnik, S., & Algina, J. (2003). Generalized eta and omega squared statistics: Measures of effect size for some common research designs. *Psychological Methods*, 8(4), 434-447.

See Also

`eta_squared()`, `anova_test()`. The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Data preparation
df <- ToothGrowth
df$dose <- as.factor(df$dose)

# Fit the model
res.aov <- aov(len ~ supp * dose, data = df)

# Effect size
omega_squared(res.aov)
partial_omega_squared(res.aov)
```

posthoc_test

Choose and Run the Appropriate Post-Hoc Test

Description

Given a one-way, independent-groups design (outcome ~ group), check the ANOVA assumptions and run the post-hoc test they imply, following the standard decision tree:

- each group normal **and** variances equal: Tukey HSD (`tukey_hsd()`);
- each group normal **but** variances unequal: Games-Howell (`games_howell_test()`);
- at least one group not normal: Dunn's test (`dunn_test()`).

Normality is assessed **per group** with the Shapiro-Wilk test applied to each group's values, routing on the smallest p-value across groups (a single non-normal group sends the data to the non-parametric test). This is deliberately not the pooled model residuals, which unequal variances would make non-normal and so hide the Games-Howell case. Homogeneity of variance is assessed with Levene's test ([levene_test\(\)](#)). Both are judged at the significance level. The function returns the chosen test's usual pairwise result, with the selected method and the assumption verdicts attached (and shown when the result is printed), so the routing is transparent rather than hidden.

To check the same assumptions before the omnibus test and read off the recommended omnibus/post-hoc pair, use [check_test_assumptions\(\)](#); its result can be passed back here through `.assumptions` so the checks are not repeated. Choosing a test by first testing its assumptions on the same data has a known cost — see the note in `?check_test_assumptions`.

See the Datanovia tutorial [Statistical Tests and Assumptions in R](#) for a worked walkthrough.

Usage

```
posthoc_test(
  data,
  formula,
  significance = 0.05,
  ...,
  .assumptions = NULL,
  omnibus = NULL
)

## S3 method for class 'posthoc_test'
print(x, ...)
```

Arguments

<code>data</code>	a data frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric outcome variable and <code>group</code> is a factor with two or more levels giving the independent groups.
<code>significance</code>	the significance level used to judge the Shapiro-Wilk and Levene assumption tests. Default is 0.05.
<code>...</code>	additional arguments forwarded to the selected post-hoc test, but only those it accepts, so an argument meant for one route does not error on another. In particular <code>p.adjust.method</code> is honoured only when Dunn's test is chosen; Tukey HSD and Games-Howell carry their own built-in adjustment and ignore it.
<code>.assumptions</code>	(optional) the tibble returned by check_test_assumptions() for the same data. When supplied its verdicts are used directly, so the Shapiro-Wilk and Levene tests are not run a second time (useful after check_test_assumptions() has already checked them). Default <code>NULL</code> (compute the assumptions here).
<code>omnibus</code>	(optional) an omnibus test result — from anova_test() , welch_anova_test() or kruskal_test() — that you already ran. If the post-hoc route chosen here belongs to a different family than that omnibus — the families being parametric equal-variance (ANOVA / Tukey), parametric unequal-variance (Welch /

Games-Howell) and non-parametric (Kruskal-Wallis / Dunn) — a warning is issued so the two stay coherent (for example an `anova_test()` omnibus followed by a Games-Howell route). Default NULL (no check).

x an object of class `posthoc_test`.

Value

the pairwise comparison table returned by the selected post-hoc test (a `tukey_hsd`, `games_howell_test` or `dunn_test` object), additionally classed `posthoc_test`. The selected method and the assumption verdicts are stored in the attributes `"posthoc.method"` and `"assumptions"`, and printed above the table.

See Also

`check_test_assumptions()`, `tukey_hsd()`, `games_howell_test()`, `dunn_test()`, `levene_test()`, `shapiro_test()`. The Datanovia tutorial: [Statistical Tests and Assumptions in R](#).

Examples

```
df <- ToothGrowth
df$dose <- as.factor(df$dose)

# Assumptions hold here, so Tukey HSD is chosen
df %>% posthoc_test(len ~ dose)
```

prop_test

Proportion Test

Description

Performs proportion tests to either evaluate the homogeneity of proportions (probabilities of success) in several groups or to test that the proportions are equal to certain given values.

Wrappers around the R base function `prop.test()` but have the advantage of performing pairwise and row-wise z-test of two proportions, the post-hoc tests following a significant chi-square test of homogeneity for 2xc and rx2 contingency tables.

See the Datanovia tutorial [Proportion Z-Test in R](#) for a worked walkthrough.

Usage

```
prop_test(
  x,
  n,
  p = NULL,
  alternative = c("two.sided", "less", "greater"),
  correct = TRUE,
  conf.level = 0.95,
  detailed = FALSE
```

```
)

pairwise_prop_test(xtab, p.adjust.method = "holm", ...)

row_wise_prop_test(xtab, p.adjust.method = "holm", detailed = FALSE, ...)
```

Arguments

<code>x</code>	a vector of counts of successes, a one-dimensional table with two entries, or a two-dimensional table (or matrix) with 2 columns, giving the counts of successes and failures, respectively.
<code>n</code>	a vector of counts of trials; ignored if <code>x</code> is a matrix or a table.
<code>p</code>	a vector of probabilities of success. The length of <code>p</code> must be the same as the number of groups specified by <code>x</code> , and its elements must be greater than 0 and less than 1.
<code>alternative</code>	a character string specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter. Only used for testing the null that a single proportion equals a given value, or that two proportions are equal; ignored otherwise.
<code>correct</code>	a logical indicating whether Yates' continuity correction should be applied where possible.
<code>conf.level</code>	confidence level of the returned confidence interval. Must be a single number between 0 and 1. Only used when testing the null that a single proportion equals a given value, or that two proportions are equal; ignored otherwise.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>xtab</code>	a cross-tabulation (or contingency table) with two columns and multiple rows (rx2 design). The columns give the counts of successes and failures respectively.
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>...</code>	Other arguments passed to the function <code>prop_test()</code> .

Value

return a data frame with some the following columns:

- `n`: the number of participants.
- `group`: the categories in the row-wise proportion tests.
- `statistic`: the value of Pearson's chi-squared test statistic.
- `df`: the degrees of freedom of the approximate chi-squared distribution of the test statistic.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the used statistical test.

- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: a vector with the sample proportions x/n .
- `estimate1`, `estimate2`: the proportion in each of the two populations.
- `alternative`: a character string describing the alternative hypothesis.
- `conf.low`, `conf.high`: Lower and upper bound on a confidence interval. a confidence interval for the true proportion if there is one group, or for the difference in proportions if there are 2 groups and `p` is not given, or `NULL` otherwise. In the cases where it is not `NULL`, the returned confidence interval has an asymptotic confidence level as specified by `conf.level`, and is appropriate to the specified alternative hypothesis.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `prop_test()`: performs one-sample and two-samples z-test of proportions. Wrapper around the function `prop.test()`.
- `pairwise_prop_test()`: pairwise comparisons between proportions, a post-hoc tests following a significant chi-square test of homogeneity for 2xc design. Wrapper around `pairwise.prop.test()`
- `row_wise_prop_test()`: performs row-wise z-test of two proportions, a post-hoc tests following a significant chi-square test of homogeneity for rx2 contingency table. The z-test of two proportions is calculated for each category (row).

See Also

The Datanovia tutorial: [Proportion Z-Test in R](#).

Examples

```
# Comparing an observed proportion to an expected proportion
#####
prop_test(x = 95, n = 160, p = 0.5, detailed = TRUE)

# Comparing two proportions
#####
# Data: frequencies of smokers between two groups
xtab <- as.table(rbind(c(490, 10), c(400, 100)))
dimnames(xtab) <- list(
  group = c("grp1", "grp2"),
  smoker = c("yes", "no")
)
xtab
# compare the proportion of smokers
prop_test(xtab, detailed = TRUE)

# Homogeneity of proportions between groups
#####
# H0: the proportion of smokers is similar in the four groups
# Ha: this proportion is different in at least one of the populations.
```

```

#
# Data preparation
grp.size <- c( 106, 113, 156, 102 )
smokers  <- c( 50, 100, 139, 80 )
no.smokers <- grp.size - smokers
xtab <- as.table(rbind(
  smokers,
  no.smokers
))
dimnames(xtab) <- list(
  Smokers = c("Yes", "No"),
  Groups = c("grp1", "grp2", "grp3", "grp4")
)
xtab

# Compare the proportions of smokers between groups
prop_test(xtab, detailed = TRUE)

# Pairwise comparison between groups
pairwise_prop_test(xtab)

# Pairwise proportion tests
#####
# Data: Titanic
xtab <- as.table(rbind(
  c(122, 167, 528, 673),
  c(203, 118, 178, 212)
))
dimnames(xtab) <- list(
  Survived = c("No", "Yes"),
  Class = c("1st", "2nd", "3rd", "Crew")
)
xtab
# Compare the proportion of survived between groups
pairwise_prop_test(xtab)

# Row-wise proportion tests
#####
# Data: Titanic
xtab <- as.table(rbind(
  c(180, 145), c(179, 106),
  c(510, 196), c(862, 23)
))
dimnames(xtab) <- list(
  Class = c("1st", "2nd", "3rd", "Crew"),
  Gender = c("Male", "Female")
)
xtab
# Compare the proportion of males and females in each category
row_wise_prop_test(xtab)

```

prop_trend_test	<i>Test for Trend in Proportions</i>
-----------------	--------------------------------------

Description

Performs chi-squared test for trend in proportion. This test is also known as Cochran-Armitage trend test.

Wrappers around the R base function `prop.trend.test()` but returns a data frame for easy data visualization.

See the Datanovia tutorial [Cochran-Armitage Trend Test in R](#) for a worked walkthrough.

Usage

```
prop_trend_test(xtab, score = NULL)
```

Arguments

xtab	a cross-tabulation (or contingency table) with two columns and multiple rows (rx2 design). The columns give the counts of successes and failures respectively.
score	group score. If NULL, the default is group number.

Value

return a data frame with some the following columns:

- n: the number of participants.
- statistic: the value of Chi-squared trend test statistic.
- df: the degrees of freedom.
- p: p-value.
- method: the used statistical test.
- p.signif: the significance level of p-values and adjusted p-values, respectively.

The **returned object has an attribute called args**, which is a list holding the test arguments.

See Also

The Datanovia tutorial: [Cochran-Armitage Trend Test in R](#).

Examples

```
# Proportion of renal stone (calculi) across age
# %%%%%%%%%%%
# Data
xtab <- as.table(rbind(
  c(384, 536, 335),
  c(951, 869, 438)
```



```

))
dimnames(xtab) <- list(
  stone = c("yes", "no"),
  age = c("30-39", "40-49", "50-59")
)
xtab
# Compare the proportion of survived between groups
prop_trend_test(xtab)

```

pull_triangle

*Pull Lower and Upper Triangular Part of a Matrix***Description**

Returns the lower or the upper triangular part of a (correlation) matrix.

Usage

```

pull_triangle(x, triangle = c("lower", "upper"), diagonal = FALSE)

pull_upper_triangle(x, diagonal = FALSE)

pull_lower_triangle(x, diagonal = FALSE)

```

Arguments

x	a (correlation) matrix
triangle	the triangle to pull. Allowed values are one of "upper" and "lower".
diagonal	logical. Default is FALSE. If TRUE, the matrix diagonal is included.

Value

an object of class `cor_mat_tri`, which is a data frame

Functions

- `pull_triangle()`: returns either the lower or upper triangular part of a matrix.
- `pull_upper_triangle()`: returns an object of class `upper_tri`, which is a data frame containing the upper triangular part of a matrix.
- `pull_lower_triangle()`: returns an object of class `lower_tri`, which is a data frame containing the lower triangular part of a matrix.

See Also

[replace_triangle\(\)](#)

Examples

```
# Data preparation
#::::::::::::::::::::::::::::::::::::
mydata <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec)
head(mydata, 3)

# Compute correlation matrix and pull triangles
#::::::::::::::::::::::::::::::::::::
# Correlation matrix
cor.mat <- cor_mat(mydata)
cor.mat

# Pull lower triangular part
cor.mat %>% pull_lower_triangle()

# Pull upper triangular part
cor.mat %>% pull_upper_triangle()
```

p_round

*Rounding and Formatting p-values***Description**

Round and format p-values. Can also mark significant p-values with stars.

Usage

```
p_round(x, ..., digits = 3)
```

```
p_format(
  x,
  ...,
  new.col = FALSE,
  digits = 2,
  accuracy = 1e-04,
  decimal.mark = ".",
  leading.zero = TRUE,
  trailing.zero = FALSE,
  add.p = FALSE,
  space = FALSE
)
```

```
p_mark_significant(
  x,
  ...,
```

```

    new.col = FALSE,
    cutpoints = c(0, 1e-04, 0.001, 0.01, 0.05, 1),
    symbols = c("****", "***", "**", "*", "")
  )

p_detect(data, type = c("all", "p", "p.adj"))

p_names()

p_adj_names()

```

Arguments

x	a numeric vector of p-values or a data frame containing a p value column. If data frame, the p-value column(s) will be automatically detected. Known p-value column names can be obtained using the functions p_names() and p_adj_names()
...	column names to manipulate in the case where x is a data frame. P value columns are automatically detected if not specified.
digits	the number of significant digits to be used.
new.col	logical, used only when x is a data frame. If TRUE, add a new column to hold the results. The new column name is created by adding, to the p column, the suffix "format" (for p_format()), "signif" (for p_mak_significant()).
accuracy	number to round to, that is the threshold value above which the function will replace the pvalue by "<0.0xxx".
decimal.mark	the character to be used to indicate the numeric decimal point.
leading.zero	logical. If FALSE, remove the leading zero.
trailing.zero	logical. If FALSE (default), remove the trailing extra zero.
add.p	logical value. If TRUE, add "p=" before the value.
space	logical. If TRUE (default) use space as separator between different elements and symbols.
cutpoints	numeric vector used for intervals
symbols	character vector, one shorter than cutpoints, used as significance symbols.
data	a data frame
type	the type of p-value to detect. Can be one of c("all", "p", "p.adj").

Value

a vector or a data frame containing the rounded/formatted p-values.

Functions

- p_round(): round p-values
- p_format(): format p-values. Add a symbol "<" for small p-values.
- p_mark_significant(): mark p-values with significance levels

- `p_detect()`: detects and returns p-value column names in a data frame.
- `p_names()`: returns known p-value column names
- `p_adj_names()`: returns known adjust p-value column names

Examples

```
# Round and format a vector of p-values
# ::::::::::::::::::::::::::::::::::::::::::::
# Format
p <- c(0.5678, 0.127, 0.045, 0.011, 0.009, 0.00002, NA)
p_format(p)

# Specify the accuracy
p_format(p, accuracy = 0.01)

# Add p and remove the leading zero
p_format(p, add.p = TRUE, leading.zero = FALSE)

# Remove space before and after "=" or "<".
p_format(p, add.p = TRUE, leading.zero = FALSE, space = FALSE)

# Mark significant p-values
# ::::::::::::::::::::::::::::::::::::::::::::
p_mark_significant(p)

# Round, the mark significant
p %>% p_round(digits = 2) %>% p_mark_significant()

# Format, then mark significant
p %>% p_format(digits = 2) %>% p_mark_significant()

# Perform stat test, format p and mark significant
# ::::::::::::::::::::::::::::::::::::::::::::
ToothGrowth %>%
  group_by(dose) %>%
  t_test(len ~ supp) %>%
  p_format(digits = 2, leading.zero = FALSE) %>%
  p_mark_significant()
```

remove_ns

Remove Non-Significant from Statistical Tests

Description

Filter out non-significant (NS) p-values from a statistical test. Can detect automatically p-value columns

Usage

```
remove_ns(stat.test, col = NULL, signif.cutoff = 0.05)
```

Arguments

stat.test	statistical test results returned by <code>rstatix</code> functions or any data frame containing a p-value column.
col	(optional) character specifying the column containing the p-value or the significance information, to be used for the filtering step. Possible values include: "p", "p.adj", "p.signif", "p.adj.signif". If missing, the function will automatically look for p.adj.signif, p.adj, p.signif, p in this order.
signif.cutoff	the significance cutoff; default is 0.05. Significance is declared at p-value <= signif.cutoff

Value

a data frame

Examples

```
# Statistical test
stat.test <- PlantGrowth %>% wilcox_test(weight ~ group)
# Remove ns: automatic detection of p-value columns
stat.test %>% remove_ns()
# Remove ns by the column p
stat.test %>% remove_ns(col = "p")
```

replace_triangle

Replace Lower and Upper Triangular Part of a Matrix

Description

Replace the lower or the upper triangular part of a (correlation) matrix.

Usage

```
replace_triangle(x, triangle = c("lower", "upper"), by = "", diagonal = FALSE)
```

```
replace_upper_triangle(x, by = "", diagonal = FALSE)
```

```
replace_lower_triangle(x, by = "", diagonal = FALSE)
```

Arguments

x	a (correlation) matrix
triangle	the triangle to replace. Allowed values are one of "upper" and "lower".
by	a replacement argument. Appropriate values are either "" or NA. Used to replace the upper, lower or the diagonal part of the matrix.
diagonal	logical. Default is FALSE. If TRUE, the matrix diagonal is included.

Value

an object of class `cor_mat_tri`, which is a data frame

Functions

- `replace_triangle()`: replaces the specified triangle by empty or NA.
- `replace_upper_triangle()`: replaces the upper triangular part of a matrix. Returns an object of class `lower_tri`.
- `replace_lower_triangle()`: replaces the lower triangular part of a matrix. Returns an object of class `lower_tri`

See Also

[pull_triangle\(\)](#)

Examples

```
# Compute correlation matrix and pull triangles
#::::::::::::::::::::::::::::::::::::::::::::::::::
# Correlation matrix
cor.mat <- mtcars %>%
  select(mpg, disp, hp, drat, wt, qsec) %>%
  cor_mat()
cor.mat

# Replace upper triangle by NA
#::::::::::::::::::::::::::::::::::::::::::::::::::
cor.mat %>% replace_upper_triangle(by = NA)

# Replace upper triangle by NA and reshape the
# correlation matrix to have unique combinations of variables
#::::::::::::::::::::::::::::::::::::::::::::::::::
cor.mat %>%
  replace_upper_triangle(by = NA) %>%
  cor_gather()
```

Description

How to use `rstatix` functions programmatically — i.e. when the variable names are held in character strings or passed into your own wrapper functions, as is common in package development, loops, and Shiny apps.

`rstatix` has two kinds of interfaces, and each supports a standard way of "programming over variables":

- **Selection interface** (functions that select columns through `...` or `vars=`, e.g. `cor_test()`, `get_summary_stats()`, `cor_mat()`, `freq_table()`). These support full tidy-evaluation: bare names, the injection operators `!!!`, the embracing operator `{ }` inside your own functions, a character vector via `vars=`, and tidyselect helpers such as `all_of()/any_of()`.
- **Formula interface** (tests that take a formula, e.g. `t_test()`, `wilcox_test()`, `kruskal_test()`, `anova_test()`, `friedman_test()`). A formula is an ordinary R object, so build it from strings with `reformulate()` or `stats::as.formula(paste(...))` and pass it in.

Details

Injecting a string directly into a raw formula (e.g. `t_test(df, y ~ {{var}})`) is **not** supported: a formula is captured as a syntax tree, not a quosure, so the embracing/injection operators do not apply there. Build the formula instead with `reformulate(rhs, lhs)` — see the examples.

In examples below, helpers that `rstatix` does not re-export are namespaced (`rlang::sym`, `dplyr::all_of`, `dplyr::across`); attach `rlang/dplyr` and you can drop the prefixes.

See Also

`cor_test()`, `get_summary_stats()`, `t_test()`.

Examples

```
# Selection interface -----

# A variable name held in a string, injected with !!
x <- "mpg"; y <- "wt"
mtcars %>% cor_test(!!rlang::sym(x), !!rlang::sym(y))

# Several names at once, spliced with !!!
vars <- c("mpg", "disp", "hp")
mtcars %>% cor_test(!!!rlang::syms(vars))

# A character vector via `vars =` (no rlang needed)
mtcars %>% cor_test(vars = c("mpg", "disp"))

# tidyselect helpers
iris %>% get_summary_stats(dplyr::all_of(c("Sepal.Length", "Sepal.Width")))
```

```

# Your own wrapper function: embrace the argument with {{ }}
my_summary <- function(data, var) {
  data %>% get_summary_stats({{ var }}, type = "mean_sd")
}
my_summary(iris, Sepal.Length)

# Formula interface -----

# Build the formula from strings with reformulate(rhs, lhs)
outcome <- "len"; group <- "supp"
ToothGrowth %>% t_test(reformulate(group, outcome))

# The same inside a wrapper function
my_test <- function(data, outcome, group) {
  data %>% t_test(reformulate(group, outcome))
}
my_test(ToothGrowth, "len", "supp")

# Programmatic grouping + a built formula (a common end-to-end pattern)
gv <- "supp"
ToothGrowth %>%
  group_by(dplyr::across(dplyr::all_of(gv))) %>%
  t_test(reformulate("dose", "len"))

```

rstatix-references	<i>References and related packages</i>
--------------------	--

Description

Where the methods in `rstatix` come from, and which other packages implement them.

Most of the statistical methods in `rstatix` are implemented in base R from the published formulas. A few call another package to compute the result, and one is adapted from another package's code; *Adapted code* below says which. Where a function records the source of its method, it does so in its own References section; cite those authors, not this package, when you report a result.

Method sources

- **Cramer's V** — Cramer, H. (1946) *Mathematical Methods of Statistics*. See `cramer_v()`.
- **Effect-size confidence intervals** (partial eta squared, Cramer's V) — obtained by inverting a noncentral distribution: Smithson, M. (2003) *Confidence Intervals*; Steiger, J. H. (2004) Beyond the F test. *Psychological Methods*, 9, 164-182. See `anova_test()` and `cramer_v()`.
- **Conover's all-pairs test** — Conover, W. J. (1999) *Practical Nonparametric Statistics*, 3rd edition. See `conover_test()` and `friedman_conover_test()`.
- **Nemenyi's all-pairs test** — Nemenyi, P. (1963) *Distribution-free Multiple Comparisons*. See `friedman_nemenyi_test()`.

- **Compact letter display** — Piepho, H.-P. (2004) An algorithm for a letter-based representation of all-pairwise comparisons. *Journal of Computational and Graphical Statistics*, 13, 456-466. See [add_cld\(\)](#).
- **Dunnett's many-to-one comparisons** — Dunnett, C. W. (1955) A multiple comparison procedure for comparing several treatments with a control. *Journal of the American Statistical Association*, 50, 1096-1121. See [dunnett_test\(\)](#).

Related packages

The following packages implement some of the same methods. `rstatix` compares its results against them while developing, and they offer functionality that `rstatix` does not:

- `effectsize` — a broad effect-size toolkit; `effectsize::cramers_v()` and `effectsize::eta_squared()` produce the same intervals as `cramer_v(ci = TRUE)`, `anova_test(ci = )`, and `eta_squared(ci = )` / `partial_eta_squared(ci = )` when called with `alternative = "two.sided"`, which is not their default.
- `DescTools` — `DescTools::CramerV()` and `DescTools::DunnettTest()`.
- `multcomp` — `multcomp::glht()` for general linear hypotheses, including Dunnett contrasts.
- `multcompView` — `multcompView::multcompLetters()` for compact letter displays.
- `PMCMRplus` — a large collection of all-pairs and many-to-one nonparametric procedures, including the Conover and Nemenyi tests.

These packages are not dependencies of `rstatix`. The values they return are recorded in the `rstatix` test suite as fixed numbers, checked against the package they came from at the time they were written; a recorded number cannot detect a later change in the package that produced it.

Adapted code

Of the functions documented in this package, [sign_test\(\)](#) is the exception to the description above. Its one- and two-sample test code, and the confidence interval it reports for the median, are adapted with modifications from `DescTools::SignTest()` and `DescTools::MedianCI()`, written by Andri Signorell. `DescTools` is distributed under GPL (>= 2); `rstatix` is distributed under GPL-2.

No other source file in `rstatix` carries a statistical method adapted from a contributed package. Argument-validation idioms are shared with base R — `check_two_samples_test_args()` follows the preamble of `stats::wilcox.test()`, as `DescTools` and `PMCMRplus` do — and those are not methods anyone cites. The test suite scans the sources for the phrases that declare an adaptation and fails when one appears in a file this section does not name; code copied without a word about it would not be caught that way.

Copying code is not the same as calling it. Several functions compute their result by calling another package, as each of them states in its own documentation: [anova_test\(\)](#) uses `car::Anova()` for type II and type III sums of squares, and `stats::aov()` for type I; [dunnett_test\(\)](#) and [emmeans_test\(\)](#) use `emmeans::emmeans()`; and [wilcox_effsize\(\)](#) takes its default statistic ($r = Z / \sqrt{N}$) from `coin::wilcoxsign_test()` or `coin::wilcox_test()`, while its `method = "rank_biserial"` alternative is computed in base R.

See Also

[cramer_v\(\)](#), [anova_test\(\)](#), [conover_test\(\)](#), [add_cld\(\)](#), [dunnett_test\(\)](#).

sample_n_by	<i>Sample n Rows By Group From a Table</i>
-------------	--

Description

sample n rows by group from a table using the [sample_n\(\)](#) function.

Usage

```
sample_n_by(data, ..., size = 1, replace = FALSE)
```

Arguments

- | | |
|---------|------------------------------|
| data | a data frame |
| ... | Variables to group by |
| size | the number of rows to select |
| replace | with or without replacement? |

Examples

```
ToothGrowth %>% sample_n_by(dose, supp, size = 2)
```

shapiro_test	<i>Shapiro-Wilk Normality Test</i>
--------------	------------------------------------

Description

Provides a pipe-friendly framework to performs Shapiro-Wilk test of normality. Support grouped data and multiple variables for multivariate normality tests. Wrapper around the R base function [shapiro.test\(\)](#). Can handle grouped data.

See the Datanovia tutorial [Normality Test in R](#) for a worked walkthrough.

Usage

```
shapiro_test(data, ..., vars = NULL)

mshapiro_test(data)
```

Arguments

data	a data frame. Columns are variables.
...	One or more unquoted expressions (or variable names) separated by commas. Used to select a variable of interest.
vars	optional character vector containing variable names. Ignored when dot vars are specified.

Value

a data frame containing the value of the Shapiro-Wilk statistic and the corresponding p.value.

Functions

- `shapiro_test()`: univariate Shapiro-Wilk normality test
- `mshapiro_test()`: multivariate Shapiro-Wilk normality test. This is a modified copy of the `mshapiro.test()` function of the package `mvnrmtest`, for internal convenience.

See Also

The Datanovia tutorial: [Normality Test in R](#).

Examples

```
# Shapiro Wilk normality test for one variable
iris %>% shapiro_test(Sepal.Length)

# Shapiro Wilk normality test for two variables
iris %>% shapiro_test(Sepal.Length, Petal.Width)

# Multivariate normality test
mshapiro_test(iris[, 1:3])
```

`sign_test`*Sign Test*

Description

Performs one-sample and two-sample sign tests.

See the Datanovia tutorial [Sign Test in R](#) for a worked walkthrough.

Usage

```

sign_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  alternative = "two.sided",
  mu = 0,
  conf.level = 0.95,
  detailed = FALSE
)

pairwise_sign_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  detailed = FALSE,
  ...
)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ treatment</code> .
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group).
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>alternative</code>	a character string specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
<code>mu</code>	a single number representing the value of the population median specified by the null hypothesis.
<code>conf.level</code>	confidence level of the interval.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>...</code>	other arguments passed to the function <code>sign_test()</code>

Value

return a data frame with some the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n, n1, n2`: Sample counts.
- `statistic`: Test statistic used to compute the p-value. That is the S-statistic (the number of positive differences between the data and the hypothesized median), with names attribute "S".
- `df, parameter`: degrees of freedom. Here, the total number of valid differences.
- `p`: p-value.
- `method`: the statistical test used to compare groups.
- `p.signif, p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: estimate of the effect size. It corresponds to the median of the differences.
- `alternative`: a character string describing the alternative hypothesis.
- `conf.low, conf.high`: Lower and upper bound on a confidence interval of the estimate.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `sign_test()`: Sign test
- `pairwise_sign_test()`: performs pairwise two sample Wilcoxon test.

Note

The sign test and the confidence interval for the median are adapted, with modifications, from `DescTools::SignTest()` and `DescTools::MedianCI()`, written by Andri Signorell. The results match `DescTools::SignTest()`.

Source

Adapted from `DescTools::SignTest()` and `DescTools::MedianCI()` (Andri Signorell), distributed under GPL (≥ 2). See [rstatix-references](#).

See Also

The Datanovia tutorial: [Sign Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# One-sample test
```

```

#::::::::::::::::::::::::::::::::::::
df %>% sign_test(len ~ 1, mu = 0)

# Two-samples paired test
#::::::::::::::::::::::::::::::::::::
df %>% sign_test(len ~ supp)

# Compare supp levels after grouping the data by "dose"
#::::::::::::::::::::::::::::::::::::
df %>%
  group_by(dose) %>%
  sign_test(data = ., len ~ supp) %>%
  adjust_pvalue(method = "bonferroni") %>%
  add_significance("p.adj")

# pairwise comparisons
#::::::::::::::::::::::::::::::::::::
# As dose contains more than two levels ==>
# pairwise test is automatically performed.
df %>% sign_test(len ~ dose)

# Comparison against reference group
#::::::::::::::::::::::::::::::::::::
# each level is compared to the ref group
df %>% sign_test(len ~ dose, ref.group = "0.5")

```

tidy.rstatix_test

Tidy an rstatix Test Result

Description

tidy() and glance() methods for objects of class rstatix_test — the result of a test function such as `t_test()`, `wilcox_test()`, `anova_test()` or `kruskal_test()`. The results are already tidy tibbles; these methods drop the internal rstatix classes and the stashed test arguments so the object passes cleanly to tools that dispatch on tidy / glance, such as broom, gtsummary and gt. Correlation results (`cor_test()`, `cor_mat()`) carry a different class and are not covered by these methods.

Usage

```

## S3 method for class 'rstatix_test'
tidy(x, ...)

## S3 method for class 'rstatix_test'
glance(x, ...)

```

Arguments

`x` an object of class `rstatix_test`, as returned by an `rstatix` test function.

`...` not used; present for compatibility with the generics.

Value

`tidy()` returns the same result as a plain tibble, one row per comparison or model term, with the internal classes and the `args` attribute removed. `glance()` returns a one-row tibble with the test method and `n`, the number of rows in the result (the number of comparisons or model terms).

Examples

```
res <- ToothGrowth %>% t_test(len ~ dose)

# A plain tibble, ready for broom / gtsummary / gt
tidy(res)

# One-row summary
glance(res)
```

tukey_hsd

Tukey Honest Significant Differences

Description

Provides a pipe-friendly framework to performs Tukey post-hoc tests. Wrapper around the function [TukeyHSD\(\)](#). It is essentially a t-test that corrects for multiple testing.

Can handle different inputs formats: `aov`, `lm`, `formula`.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
tukey_hsd(x, ...)
```

Default S3 method:

```
tukey_hsd(x, ...)
```

S3 method for class 'lm'

```
tukey_hsd(x, ...)
```

S3 method for class 'data.frame'

```
tukey_hsd(x, formula, ...)
```

Arguments

<code>x</code>	an object of class <code>aov</code> , <code>lm</code> or <code>data.frame</code> containing the variables used in the formula.
<code>...</code>	other arguments passed to the function <code>TukeyHSD()</code> . These include: • which: A character vector listing terms in the fitted model for which the intervals should be calculated. Defaults to all the terms. • ordered: A logical value indicating if the levels of the factor should be ordered according to increasing average in the sample before taking differences. If <code>ordered</code> is <code>true</code> then the calculated differences in the means will all be positive. The significant differences will be those for which the lwr end point is positive.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>data</code>	a <code>data.frame</code> containing the variables in the formula.

Value

a tibble data frame containing the results of the different comparisons.

Methods (by class)

- `tukey_hsd(default)`: performs tukey post-hoc test from `aov()` results.
- `tukey_hsd(lm)`: performs tukey post-hoc test from `lm()` model.
- `tukey_hsd(data.frame)`: performs tukey post-hoc tests using `data` and `formula` as inputs. ANOVA will be automatically performed using the function `aov()`

See Also

The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Data preparation
df <- ToothGrowth
df$dose <- as.factor(df$dose)
# Tukey HSD from ANOVA results
aov(len ~ dose, data = df) %>% tukey_hsd()

# two-way anova with interaction
aov(len ~ dose*supp, data = df) %>% tukey_hsd()

# Tukey HSD from lm() results
lm(len ~ dose, data = df) %>% tukey_hsd()

# Tukey HSD from data frame and formula
tukey_hsd(df, len ~ dose)
```



```
# Tukey HSD using grouped data
df %>%
  group_by(supp) %>%
  tukey_hsd(len ~ dose)
```

t_test

T-test

Description

Provides a pipe-friendly framework to performs one and two sample t-tests.

See the Datanovia tutorial [T-Test in R](#) for a worked walkthrough.

Usage

```
t_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  paired = FALSE,
  var.equal = FALSE,
  alternative = "two.sided",
  mu = 0,
  conf.level = 0.95,
  detailed = FALSE,
  id = NULL,
  error.as.na = FALSE,
  effect.size = FALSE
)
```

```
pairwise_t_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  paired = FALSE,
  pool.sd = !paired,
  detailed = FALSE,
  ...,
  effect.size = FALSE
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>paired</code>	a logical indicating whether you want a paired test.
<code>var.equal</code>	a logical variable indicating whether to treat the two variances as being equal. If TRUE then the pooled variance is used to estimate the variance otherwise the Welch (or Satterthwaite) approximation to the degrees of freedom is used.
<code>alternative</code>	a character string specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
<code>mu</code>	a number specifying an optional parameter used to form the null hypothesis.
<code>conf.level</code>	confidence level of the interval.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>id</code>	(optional) character string specifying the column that contains the sample/subject identifier, used only for a paired test (<code>paired = TRUE</code>). When supplied, observations of the two compared groups are matched by <code>id</code> (instead of by row order), and only subjects present in both groups are used. For more than two groups, the matching is done independently for each pairwise comparison, so different comparisons can be based on different numbers of pairs (per-comparison pairwise deletion). This makes paired tests work when some observations are missing or the groups have unequal sizes. The default (<code>id = NULL</code>) keeps the previous behaviour (groups paired in row order).
<code>error.as.na</code>	logical. If TRUE, a comparison that cannot be computed (for example a group with fewer than two observations, or data that are essentially constant) returns an NA result row with a warning instead of stopping with an error; the other comparisons (or groups, for a grouped analysis) are still computed. Default is FALSE (the comparison errors as before).
<code>effect.size</code>	logical. Default is FALSE. If TRUE, a <code>cohens.d</code> column (Cohen's d) and its magnitude are added. <code>t_test()</code> reports the per-pair d (consistent with its per-pair t-test); <code>pairwise_t_test()</code> with the default <code>pool.sd = TRUE</code> reports the

	common-SD (pooled-model) d so the estimate and the pooled-SD p-value share a variance basis. For a paired test the d is computed on the same paired differences the test used, matched by id when it is supplied.
pool.sd	logical value used in the function <code>pairwise_t_test()</code> . Switch to allow/disallow the use of a pooled SD. The <code>pool.sd = TRUE</code> (default) calculates a common SD for all groups and uses that for all comparisons (this can be useful if some groups are small). This method does not actually call <code>t.test</code> , so extra arguments are ignored. Pooling does not generalize to paired tests so <code>pool.sd</code> and <code>paired</code> cannot both be <code>TRUE</code> . If <code>pool.sd = FALSE</code> the standard two sample t-test is applied to all possible pairs of groups. This method calls the <code>t.test()</code> , so extra arguments, such as <code>var.equal</code> are accepted.
...	other arguments to be passed to the function <code>t.test</code> .

Details

- If a list of comparisons is specified, the result of the pairwise tests is filtered to keep only the comparisons of interest. The p-value is adjusted after filtering.
- For a grouped data, if pairwise test is performed, then the p-values are adjusted for each group level independently.

Value

return a data frame with some the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n, n1, n2`: Sample counts.
- `statistic`: Test statistic used to compute the p-value.
- `df`: degrees of freedom.
- `p`: p-value.
- `p.adj`: the adjusted p-value.
- `method`: the statistical test used to compare groups.
- `p.signif, p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: estimate of the effect size. It corresponds to the estimated mean or difference in means depending on whether it was a one-sample test or a two-sample test. For a two-sample test the difference is taken as `estimate1 - estimate2`, i.e. `mean(group1) - mean(group2)` (following `t.test`).
- `estimate1, estimate2`: show the mean values of the two groups, respectively, for independent samples t-tests.
- `alternative`: a character string describing the alternative hypothesis.
- `conf.low, conf.high`: Lower and upper bound on a confidence interval.

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `t_test()`: t test
- `pairwise_t_test()`: performs pairwise two sample t-test. Wrapper around the R base function `pairwise.t.test`.

Note

On the sign of statistic and estimate when a `ref.group` is specified: the reference group is taken as `group1` and the other group as `group2`, and the difference is computed as `estimate = mean(group1) - mean(group2) = mean(ref.group) - mean(other)` (the `t.test` convention). A positive statistic/estimate therefore means the value is higher in the reference group. To orient results so that a positive sign means "higher in the non-reference group", flip the sign yourself, e.g. `mutate(statistic = -statistic, estimate = -estimate)`.

See Also

[rstatix-programming](#) for building the formula from variable names held in strings (e.g. `reformulate()`).
The Datanovia tutorial: [T-Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# One-sample test
#::::::::::::::::::::::::::::::::::::
df %>% t_test(len ~ 1, mu = 0)

# Two-samples unpaired test
#::::::::::::::::::::::::::::::::::::
df %>% t_test(len ~ supp)

# Two-samples paired test
#::::::::::::::::::::::::::::::::::::
df %>% t_test (len ~ supp, paired = TRUE)

# Compare supp levels after grouping the data by "dose"
#::::::::::::::::::::::::::::::::::::
df %>%
  group_by(dose) %>%
  t_test(data = ., len ~ supp) %>%
  adjust_pvalue(method = "bonferroni") %>%
  add_significance("p.adj")

# pairwise comparisons
#::::::::::::::::::::::::::::::::::::
# As dose contains more than two levels ==>
# pairwise test is automatically performed.
```

```
df %>% t_test(len ~ dose)

# Comparison against reference group
#::::::::::::::::::::::::::::::::::::
# each level is compared to the ref group
df %>% t_test(len ~ dose, ref.group = "0.5")

# Comparison against all
#::::::::::::::::::::::::::::::::::::
df %>% t_test(len ~ dose, ref.group = "all")
```

welch_anova_test	<i>Welch One-Way ANOVA Test</i>
------------------	---------------------------------

Description

Tests for equal means in a one-way design (not assuming equal variance). A wrapper around the base function `oneway.test()`. This is an alternative to the standard one-way ANOVA in the situation where the homogeneity of variance assumption is violated.

See the Datanovia tutorial [One-Way ANOVA in R](#) for a worked walkthrough.

Usage

```
welch_anova_test(data, formula)
```

Arguments

data	a data frame containing the variables in the formula.
formula	a formula specifying the ANOVA model similar to <code>aov</code> . Can be of the form <code>y ~ group</code> where <code>y</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .

Value

return a data frame with the following columns:

- `.y.`: the y variable used in the test.
- `n`: sample count.
- `statistic`: the value of the test statistic.
- `p`: p-value.
- `method`: the statistical test used to compare groups.

See Also

The Datanovia tutorial: [One-Way ANOVA in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth
df$dose <- as.factor(df$dose)

# Welch one-way ANOVA test (not assuming equal variance)
#::::::::::::::::::::::::::::::::::::
df %>% welch_anova_test(len ~ dose)

# Grouped data
#::::::::::::::::::::::::::::::::::::
df %>%
  group_by(supp) %>%
  welch_anova_test(len ~ dose)
```

wilcox_effsize

Wilcoxon Effect Size

Description

Compute Wilcoxon effect size (r) for:

- one-sample test (Wilcoxon one-sample signed-rank test);
- paired two-samples test (Wilcoxon two-sample paired signed-rank test) and
- independent two-samples test (Mann-Whitney, two-sample rank-sum test).

It can also return confidence intervals by bootstrap.

The effect size r is calculated as Z statistic divided by square root of the sample size (N) ($Z/\sqrt{N}$). The Z value is extracted from either `coin::wilcoxsign_test()` (case of one- or paired-samples test) or `coin::wilcox_test()` (case of independent two-samples test).

Here, N is the number of independent observations contributing to the test: the total sample size for the independent two-samples test, and the **number of pairs** (equivalently, the number of difference scores) for the one-sample and paired tests. This is because the paired test reduces to a one-sample signed-rank test on the pairwise differences, so each pair counts once. This convention matches the default of `rcompanion::wilcoxonPairedR()` (its `cases = TRUE` setting).

Some references instead define N as the total number of observations, i.e. twice the number of pairs (Field, 2012; Tomczak & Tomczak, 2014), which yields a smaller r . If you need that convention for a paired test, divide the reported r (or the Z) by $\sqrt{2}$; it is also available via `rcompanion::wilcoxonPairedR(..., cases = FALSE)`.

The r value varies from 0 to close to 1. The interpretation values for r commonly in published literature and on the internet are: $0.10 - < 0.3$ (small effect), $0.30 - < 0.5$ (moderate effect) and ≥ 0.5 (large effect).

See the Datanovia tutorial [Wilcoxon Test in R](#) for a worked walkthrough.

Usage

```
wilcox_effsize(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  paired = FALSE,
  alternative = "two.sided",
  mu = 0,
  ci = FALSE,
  conf.level = 0.95,
  ci.type = "perc",
  nboot = 1000,
  detailed = FALSE,
  ...,
  boot.parallel = getOption("boot.parallel", "no"),
  boot.ncpus = getOption("boot.ncpus", 1L),
  method = c("r", "rank_biserial")
)
```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
<code>paired</code>	a logical indicating whether you want a paired test.
<code>alternative</code>	a character string specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
<code>mu</code>	a number specifying an optional parameter used to form the null hypothesis.
<code>ci</code>	If TRUE, returns confidence intervals by bootstrap. May be slow.
<code>conf.level</code>	The level for the confidence interval.
<code>ci.type</code>	The type of confidence interval to use. Can be any of "norm", "basic", "perc", or "bca". Passed to <code>boot::boot.ci</code> .
<code>nboot</code>	The number of replications to use for bootstrap.

detailed	logical value. Default is FALSE. If TRUE, and method = "r", the output additionally includes the Z statistic (extracted from the coin package and used to compute $r = Z/\sqrt{N}$), the p-value (p) and the test method/alternative, so the effect size and the underlying Z are reported together in one data frame. The rank-biserial metric (method = "rank_biserial") has no underlying Z, so those extra columns are not meaningful for it.
...	Additional arguments passed to the functions <code>coin::wilcoxsign_test()</code> (case of one- or paired-samples test) or <code>coin::wilcox_test()</code> (case of independent two-samples test).
boot.parallel	The type of parallel operation to be used when computing the bootstrap confidence interval. Allowed values are "no" (default), "multicore" and "snow". Passed to <code>boot()</code> . Defaults to <code>getOption("boot.parallel", "no")</code> , so it can also be set globally with <code>options(boot.parallel = "multicore")</code> . Only used when <code>ci = TRUE</code> .
boot.ncpus	Integer. The number of processes to be used in the parallel bootstrap. Defaults to <code>getOption("boot.ncpus", 1L)</code> . Note that <code>boot.parallel</code> has no effect unless <code>boot.ncpus > 1</code> . Only used when <code>ci = TRUE</code> .
method	the effect-size metric. Either "r" (default) for the rank correlation $r = Z / \sqrt{N}$, or "rank_biserial" for the rank-biserial correlation — Cliff's delta for an independent-samples test (equal to <code>cliff_delta()</code>) or the matched-pairs rank-biserial for a paired test, both equal to <code>effectsize::rank_biserial()</code> . The independent-samples case is labelled with the Romano et al. magnitude thresholds (those thresholds define Cliff's delta); the paired case carries no magnitude (NA), because no threshold set is calibrated for the matched-pairs rank-biserial. The confidence interval (<code>ci = TRUE</code>) is a percentile bootstrap.

Value

return a data frame with some of the following columns:

- `.y.`: the y variable used in the test.
- `group1, group2`: the compared groups in the pairwise tests.
- `n, n1, n2`: Sample counts.
- `effsize`: estimate of the effect size (r value).
- `magnitude`: magnitude of effect size.
- `conf.low, conf.high`: lower and upper bound of the effect size confidence interval.
- `statistic`: the Z statistic and p: the p-value (only when `detailed = TRUE`).

References

Maciej Tomczak and Ewa Tomczak. The need to report effect size estimates revisited. An overview of some recommended measures of effect size. *Trends in Sport Sciences*. 2014; 1(21):19-25.

See Also

The Datanovia tutorial: [Wilcoxon Test in R](#).

Examples

```

if(require("coin")){

# One-sample Wilcoxon test effect size
ToothGrowth %>% wilcox_effsize(len ~ 1, mu = 0)

# Independent two-samples wilcoxon effect size
ToothGrowth %>% wilcox_effsize(len ~ supp)

# Paired-samples wilcoxon effect size
ToothGrowth %>% wilcox_effsize(len ~ supp, paired = TRUE)

# Pairwise comparisons
ToothGrowth %>% wilcox_effsize(len ~ dose)

# Grouped data
ToothGrowth %>%
  group_by(supp) %>%
  wilcox_effsize(len ~ dose)

}

```

wilcox_test

Wilcoxon Tests

Description

Provides a pipe-friendly framework to performs one and two sample Wilcoxon tests.

See the Datanovia tutorial [Wilcoxon Test in R](#) for a worked walkthrough.

Usage

```

wilcox_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  paired = FALSE,
  exact = NULL,
  alternative = "two.sided",
  mu = 0,
  conf.level = 0.95,
  detailed = FALSE,
  id = NULL,
  error.as.na = FALSE,
  effect.size = FALSE

```

```

)

pairwise_wilcox_test(
  data,
  formula,
  comparisons = NULL,
  ref.group = NULL,
  p.adjust.method = "holm",
  detailed = FALSE,
  ...,
  effect.size = FALSE
)

```

Arguments

<code>data</code>	a data.frame containing the variables in the formula.
<code>formula</code>	a formula of the form <code>x ~ group</code> where <code>x</code> is a numeric variable giving the data values and <code>group</code> is a factor with one or multiple levels giving the corresponding groups. For example, <code>formula = TP53 ~ cancer_group</code> .
<code>comparisons</code>	A list of length-2 vectors specifying the groups of interest to be compared. For example to compare groups "A" vs "B" and "B" vs "C", the argument is as follow: <code>comparisons = list(c("A", "B"), c("B", "C"))</code>
<code>ref.group</code>	a character string specifying the reference group. If specified, for a given grouping variable, each of the group levels will be compared to the reference group (i.e. control group). If <code>ref.group = "all"</code> , pairwise two sample tests are performed for comparing each grouping variable levels against all (i.e. basemean).
<code>p.adjust.method</code>	method to adjust p values for multiple comparisons. Used when pairwise comparisons are performed. Allowed values include "holm", "hochberg", "hommel", "bonferroni", "BH", "BY", "fdr", "none". If you don't want to adjust the p value (not recommended), use <code>p.adjust.method = "none"</code> .
<code>paired</code>	a logical indicating whether you want a paired test.
<code>exact</code>	a logical indicating whether an exact p-value should be computed.
<code>alternative</code>	a character string specifying the alternative hypothesis, must be one of "two.sided" (default), "greater" or "less". You can specify just the initial letter.
<code>mu</code>	a number specifying an optional parameter used to form the null hypothesis.
<code>conf.level</code>	confidence level of the interval.
<code>detailed</code>	logical value. Default is FALSE. If TRUE, a detailed result is shown.
<code>id</code>	(optional) character string specifying the column that contains the sample/subject identifier, used only for a paired test (<code>paired = TRUE</code>). When supplied, observations of the two compared groups are matched by <code>id</code> (instead of by row order), and only subjects present in both groups are used. For more than two groups, the matching is done independently for each pairwise comparison, so different comparisons can be based on different numbers of pairs (per-comparison pairwise

	deletion). This makes paired tests work when some observations are missing or the groups have unequal sizes. The default (<code>id = NULL</code>) keeps the previous behaviour (groups paired in row order).
<code>error.as.na</code>	logical. If TRUE, a comparison that cannot be computed (for example a group with fewer than two observations, or data that are essentially constant) returns an NA result row with a warning instead of stopping with an error; the other comparisons (or groups, for a grouped analysis) are still computed. Default is FALSE (the comparison errors as before).
<code>effect.size</code>	logical. Default is FALSE. If TRUE, a rank effect-size column is added: for an independent-samples test <code>cliff.delta</code> (Cliff's delta) and its magnitude; for a paired test the matched-pairs <code>rank.biserial</code> correlation (no magnitude, as no threshold set is calibrated for it), computed on the same paired differences the test used (matched by <code>id</code> when supplied). Not defined for a one-sample test.
<code>...</code>	other arguments to be passed to the function <code>wilcox.test</code> .

Details

- `pairwise_wilcox_test()` applies the standard two sample Wilcoxon test to all possible pairs of groups. This method calls the `wilcox.test()`, so extra arguments are accepted.
- If a list of comparisons is specified, the result of the pairwise tests is filtered to keep only the comparisons of interest. The p-value is adjusted after filtering.
- For a grouped data, if pairwise test is performed, then the p-values are adjusted for each group level independently.
- a nonparametric confidence interval and an estimator for the pseudomedian (one-sample case) or for the difference of the location parameters $x - y$ is computed, where x and y are the compared samples or groups. The column `estimate` and the confidence intervals are displayed in the test result when the option `detailed = TRUE` is specified in the `wilcox_test()` and `pairwise_wilcox_test()` functions. Read more about the calculation of the estimate in the details section of the R base function `wilcox.test()` documentation by typing `?wilcox.test` in the R console.
- With `effect.size = TRUE`, an independent-samples test is annotated with Cliff's delta and a paired test with the matched-pairs rank-biserial correlation, $(R^+ - R^-)/(R^+ + R^-)$ over the signed ranks of the paired differences (Kerby, 2014).

Value

return a data frame with some of the following columns:

- `.y.:` the y variable used in the test.
- `group1, group2:` the compared groups in the pairwise tests.
- `n, n1, n2:` Sample counts.
- `statistic:` Test statistic used to compute the p-value.
- `p:` p-value.
- `p.adj:` the adjusted p-value.
- `method:` the statistical test used to compare groups.

- `p.signif`, `p.adj.signif`: the significance level of p-values and adjusted p-values, respectively.
- `estimate`: an estimate of the location parameter (Only present if argument `detailed = TRUE`). This corresponds to the pseudomedian (for one-sample case) or to the difference of the location parameter (for two-samples case).
 - The pseudomedian of a distribution F is the median of the distribution of $(u+v)/2$, where u and v are independent, each with distribution F . If F is symmetric, then the pseudomedian and median coincide.
 - Note that in the two-sample case the estimator for the difference in location parameters does not estimate the difference in medians (a common misconception) but rather the median of the difference between a sample from x and a sample from y .
- `conf.low`, `conf.high`: a confidence interval for the location parameter. (Only present if argument `conf.int = TRUE`.)

The **returned object has an attribute called `args`**, which is a list holding the test arguments.

Functions

- `wilcox_test()`: Wilcoxon test
- `pairwise_wilcox_test()`: performs pairwise two sample Wilcoxon test.

Note

When a `ref.group` is specified, the reference group is taken as `group1` and the other group as `group2`, and the comparison is computed as `group1` versus `group2` (i.e. `ref.group` versus the other group), following the `wilcox.test` convention. With `detailed = TRUE`, the estimate is the Hodges-Lehmann location shift of `group1` relative to `group2`, so a positive estimate means the reference group is shifted higher; flip its sign (`mutate(estimate = -estimate)`) if you want a positive sign to mean "higher in the non-reference group". (The statistic is the rank-sum/signed-rank W , which is not a signed difference.)

References

Kerby, D. S. (2014). The simple difference formula: An approach to teaching nonparametric correlation. *Comprehensive Psychology*, 3, 11.IT.3.1.

See Also

The Datanovia tutorial: [Wilcoxon Test in R](#).

Examples

```
# Load data
#::::::::::::::::::::::::::::::::::::
data("ToothGrowth")
df <- ToothGrowth

# One-sample test
#::::::::::::::::::::::::::::::::::::
```

```

df %>% wilcox_test(len ~ 1, mu = 0)

# Two-samples unpaired test
#::::::::::::::::::::::::::::::::::::
df %>% wilcox_test(len ~ supp)

# Two-samples paired test
#::::::::::::::::::::::::::::::::::::
df %>% wilcox_test (len ~ supp, paired = TRUE)

# Compare supp levels after grouping the data by "dose"
#::::::::::::::::::::::::::::::::::::
df %>%
  group_by(dose) %>%
  wilcox_test(data = ., len ~ supp) %>%
  adjust_pvalue(method = "bonferroni") %>%
  add_significance("p.adj")

# pairwise comparisons
#::::::::::::::::::::::::::::::::::::
# As dose contains more than two levels ==>
# pairwise test is automatically performed.
df %>% wilcox_test(len ~ dose)

# Comparison against reference group
#::::::::::::::::::::::::::::::::::::
# each level is compared to the ref group
df %>% wilcox_test(len ~ dose, ref.group = "0.5")

# Comparison against all
#::::::::::::::::::::::::::::::::::::
df %>% wilcox_test(len ~ dose, ref.group = "all")

```

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