

# Package ‘rworkflows’

July 15, 2023

**Type** Package

**Title** Test, Document, Containerise, and Deploy R Packages

**Version** 0.99.11

**Description** Reproducibility is essential to the progress of research, yet achieving it remains elusive even in computational fields. Continuous Integration (CI) platforms offer a powerful way to launch automated workflows to check and document code, but often require considerable time, effort, and technical expertise to setup. We therefore developed the rworkflows suite to make robust CI workflows easy and freely accessible to all R package developers. rworkflows consists of 1) a CRAN/Bioconductor-compatible R package template, 2) an R package to quickly implement a standardised workflow, and 3) a centrally maintained GitHub Action.

**URL** <https://github.com/neurogenomics/rworkflows>, <https://CRAN.R-project.org/package=rworkflows>

**BugReports** <https://github.com/neurogenomics/rworkflows/issues>

**Encoding** UTF-8

**biocViews** WorkflowManagement

**Depends** R (>= 4.1)

**Imports** stats,  
here,  
yaml,  
utils,  
desc,  
badger,  
renv,  
biocViews,  
tools,  
methods,  
BiocPkgTools,  
BiocManager,  
data.table

**Suggests** markdown,  
rmarkdown,  
remotes,  
knitr,  
BiocStyle,

covr,  
testthat (>= 3.0.0),  
htmltools

**RoxygenNote** 7.2.3

**VignetteBuilder** knitr

**License** GPL-3

**Config/testthat/edition** 3

**R topics documented:**

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bioc_r_versions	<i>Bioconductor / R versions</i>
-----------------	----------------------------------

---

**Description**

Get the respective version of R for a given version of **Bioconductor**.

**Usage**

bioc\_r\_versions(bioc\_version = NULL, depth = NULL)

**Arguments**

- bioc\_version    Version of Bioc to return info for. Can be:
- "devel"Get the current development version of Bioc.
  - "release"Get the current release version of Bioc.
  - <numeric>A specific Bioc version number (e.g. 3.16).
  - NULLReturn info for all Bioc versions.

- depth                      How many levels deep into the R version to include. For example, is the R version number is "4.2.0", the following depths would return:
- depth=NULL: "4.2.0"
  - depth=1: "4"
  - depth=2: "4.2"
  - depth=3: "4.2.0"

### Value

Named list of Bioc/R versions

### Examples

```
ver <- bioc_r_versions(bioc_version="devel")
```

---

construct_authors	<i>Construct authors</i>
-------------------	--------------------------

---

### Description

Helper function to construct an author list for a *DESCRIPTION* file. Returns a template when authors is not provided (default).

### Usage

```
construct_authors(
  authors = NULL,
  template = c(utils::person(given = "yourGivenName", family = "yourFamilyName", role =
    c("cre"), email = "yourEmail@email.com", comment = c(ORCID = "yourOrcidId")))
)
```

### Arguments

- authors                      A list of authors who contributed to your R package, each provided as objects of class [person](#). By default, if an Authors field already exists in the *DESCRIPTION* file, the original values are kept. Otherwise, a template [person](#) list is created using the [construct\\_authors](#).
- template                      Default value to return when authors=NULL.

### Value

Named list in [person](#) format.

### Examples

```
authors <- construct_authors()
```

---

construct_runners	<i>Construct runners</i>
-------------------	--------------------------

---

## Description

Construct runner configurations across multiple Operating Systems (OS) for GitHub Actions workflow.

## Usage

```
construct_runners(
  os = c("ubuntu-latest", "macOS-latest", "windows-latest"),
  bioc = stats::setNames(list("devel", "release", "release"), os),
  r = stats::setNames(list("auto", "auto", "auto"), os),
  cont = stats::setNames(list(paste0("bioconductor/bioconductor_docker:", bioc[[1]]),
    NULL, NULL), os),
  rspm = stats::setNames(list(paste0("https://packagemanager.rstudio.com/",
    "cran/__linux__/focal/release"), NULL, NULL), os),
  versions_explicit = FALSE
)
```

## Arguments

os	Which OS to launch GitHub Actions on.
bioc	Which Bioconductor version to use on each OS. See <a href="#">bioc_r_versions</a> documentation for all options.
r	Which R version to use on each OS.
cont	Which Docker container to use on each OS (NULL means no container will be used for that OS). See <a href="#">here</a> for a list of all official Bioconductor Docker container versions.
rspm	Which R repository manager to use on each OS (NULL means the default will be used for that OS).
versions_explicit	Specify R/Bioc versions explicitly (e.g. r: 4.2.0, bioc: 3.16) as opposed to flexibly (e.g. r: "latest", bioc: "release").

## Value

Named list of configurations for each runner OS.

## Examples

```
runners <- construct_runners()
```

---

dt_to_desc	<i>data.table</i> to <i>desc</i>
------------	----------------------------------

---

## Description

Convert [data.table](#) containing the parsed *DESCRIPTION* file data and convert each of them to [desc](#) format.

## Usage

```
dt_to_desc(db, refs = NULL, verbose = TRUE)
```

## Arguments

db	A <a href="#">data.table</a> where each row is a different R package and each column is a field from the <i>DESCRIPTION</i> file.
refs	Reference for one or more GitHub repository in owner/repo format (e.g. "neurogenomics/rworkflows") or an R package name (e.g. "rworkflows").
verbose	Print messages.

## Value

A named list of [desc](#) objects.

## Examples

```
db <- BiocPkgTools::biocPkgList()
dl <- dt_to_desc(db=db, refs="GenomicRanges")
```

---

fill_description	<i>Fill</i> DESCRIPTION
------------------	-------------------------

---

## Description

Fill out a *DESCRIPTION* file, such as (but not limited to) the one provided by the [templateR](#) R package template. For any given field, set its corresponding argument as follows to get certain behaviour:

- NULL: Keeps the current value.
- NA: Removes the field from the *DESCRIPTION* file entirely.

## Usage

```
fill_description(
  path = here::here("DESCRIPTION"),
  package,
  title,
  description,
  github_owner = NULL,
  github_repo = package,
  authors = construct_authors(authors = NULL),
  depends = paste0("R ", "(>= ", bioc_r_versions(bioc_version = "devel", depth = 2)$r,
    ")"),
  imports = infer_deps(which = "Imports", add_newlines = TRUE),
  suggests = infer_deps(which = "Suggests", add_newlines = TRUE),
  remotes = NULL,
  version = NULL,
  license = NULL,
  encoding = NULL,
  vignettebuilder = NULL,
  biocviews = infer_biocviews(pkgdir = dirname(path), add_newlines = TRUE),
  url = paste0("https://github.com/", github_owner, "/", github_repo),
  bugreports = paste0(url, "/issues"),
  save_path = path,
  verbose = TRUE,
  fields = list()
)
```

## Arguments

path	Path to the <i>DESCRIPTION</i> file.
package	The name of your R package.
title	The title of your R package.
description	The description of your R package.
github_owner	The owner of your R package's GitHub repository. Can be inferred from the URL field in the <i>DESCRIPTION</i> file if this has already been filled out.
github_repo	The name of your R package's GitHub repository.
authors	A list of authors who contributed to your R package, each provided as objects of class <a href="#">person</a> . By default, if an Authors field already exists in the <i>DESCRIPTION</i> file, the original values are kept. Otherwise, a template <a href="#">person</a> list is created using the <a href="#">construct_authors</a> .
depends	R package Depends. Defaults to the version of R that the current development version of Bioconductor depends on.
imports	R package Imports. These dependencies will be automatically installed with your R package.
suggests	R package Suggests. These dependencies will NOT be automatically installed with your R package, unless otherwise specified by users during installation
remotes	R package Remotes
version	The current version of your R package (e.g 0.99.0).
license	R package license. See <a href="#">here for guidance</a> .

encoding	R package Encoding.
vignettebuilder	R package VignetteBuilder.
biocviews	Standardised <b>biocViews</b> terms used to describe your package. Defaults to automatically recommending terms using the <a href="#">infer_biocviews</a> function. Note that non-Bioconductor packages (e.g. CRAN) can also use this field.
url	URL where your R package is distributed from (e.g. GitHub repository, Bioconductor page, and/or CRAN page). Can be a single character string or a character vector.
bugreports	A URL where users of your package should go if they encounter bugs or have feature requests.
save_path	Path to save the updated <i>DESCRIPTION</i> file to. Defaults to overwriting the input file (path). Set to NULL if you wish to only return the <a href="#">description</a> object without writing to any file.
verbose	Print messages.
fields	A named list of additional fields to fill the <i>DESCRIPTION</i> file with: e.g. <code>list(RoxygenNote=7.2.3)</code>

### Value

An object of class [description](#).

### Examples

```
#### Get example DESCRIPTION file ####
url <- "https://github.com/neurogenomics/templateR/raw/master/DESCRIPTION"
path <- tempfile(fileext = "DESCRIPTION")
utils::download.file(url,path)

#### Fill out DESCRIPTION file ####
d <- fill_description(
  path = path,
  package = "MyPackageName",
  title = "This Package Does Awesome Stuff",
  description = paste(
    "MyPackageName does several awesome things.",
    "Describe thing1.",
    "Describe thing2.",
    "Describe thing3."
  ),
  github_owner = "OwnerName",
  biocviews = c("Genetics", "SystemsBiology"))
```

---

get\_description

*Get DESCRIPTION*


---

### Description

The **Liam Neeson** of *DESCRIPTION* file functions.

1. I will look for you,
2. I will find you,

3. —and I will import you into a neatly parsed R object.

Uses a variety of alternative methods, including searching locally and on GitHub (whenever possible). Prioritises the fastest methods that do not involve downloading files first.

### Usage

```
get_description(
  refs = NULL,
  paths = here::here("DESCRIPTION"),
  cache_dir = tools::R_user_dir(package = "rworkflows", which = "cache"),
  force_new = FALSE,
  use_wd = TRUE,
  use_repos = TRUE,
  verbose = TRUE
)
```

### Arguments

refs	Reference for one or more GitHub repository in owner/repo format (e.g. "neurogenomics/rworkflows") or an R package name (e.g. "rworkflows").
paths	Paths to <i>DESCRIPTION</i> file(s) R package(s).
cache_dir	Directory where to cache downloaded files.
force_new	Ignore cached files and re-download them instead.
use_wd	Search the local working directory (and the one above it) for <i>DESCRIPTION</i> files.
use_repos	Use R standard R package repositories like CRAN and Bioc to find <i>DESCRIPTION</i> files.
verbose	Print messages.

### Value

A named list of packageDescription objects.

### Examples

```
d <- get_description(refs="neurogenomics/rworkflows")
```

---

get\_hex

*Get hex*


---

### Description

Get the URL of a hex sticker for a given R package (if one exists).



**Usage**

```
get_hex(
  refs = NULL,
  paths = here::here("DESCRIPTION"),
  hex_path = "inst/hex/hex.png",
  branch = c("master", "main", "dev"),
  hex_height = 300,
  check_url = TRUE,
  add_html = TRUE,
  verbose = TRUE
)
```

**Arguments**

refs	Reference for one or more GitHub repository in owner/repo format (e.g. "neurogenomics/rworkflows" or an R package name (e.g. "rworkflows")).
paths	Paths to <i>DESCRIPTION</i> file(s) R package(s).
hex_path	Path to hex sticker file.
branch	Name of the GitHub repository branch to use.
hex_height	Height of the hex sticker in pixels (when add_hex=TRUE).
check_url	Check whether the URL actually exists.
add_html	Wrap the URL in an html "img" tag and set its height with hex_height.
verbose	Print messages.

**Value**

URL

**Examples**

```
hex_url <- get_hex(refs=c("neurogenomics/rworkflows",
                          "neurogenomics/echolocator"))
```

---

infer\_biocviews

*Infer biocViews*


---

**Description**

Infer the best terms to fill the biocViews field in your *DESCRIPTION* file based on the code within your R package. By default, also includes any biocViews that are already present in the *DESCRIPTION* file. Please see the [Bioconductor website](#) for more details.

## Usage

```
infer_biocviews(
  pkgdir = here::here(),
  branch = c("Software", "AnnotationData", "ExperimentData")[1],
  type = c("recommended", "current", "remove"),
  keep_current = TRUE,
  include_branch = TRUE,
  biocviews = NULL,
  add_newlines = FALSE,
  verbose = TRUE
)
```

## Arguments

pkgdir	The path of the package Directory.
branch	The branch which your package will belong to. It can be either 'Software', 'AnnotationData' or 'ExperimentData'.
type	Which element of the <a href="#">recommendBiocViews</a> results list to return. If a vector is supplied, only the first value will be used.
keep_current	Keep any biocViews terms that are already included in the <i>DESCRIPTION</i> file.
include_branch	Whether to include the branch argument as one of the returned biocViews.
biocviews	User-supplied biocViews terms to include in addition to the automated recommendations.
add_newlines	Prefix each package name with a newline character and two spaces. This is useful for formatting <i>DESCRIPTION</i> files.
verbose	Print messages.

## Examples

```
## Don't run simply bc biocViews::recommendBiocViews is unable
## to find the DESCRIPTION file when running examples.
## Not run:
biocviews <- infer_biocviews()

## End(Not run)
```

---

infer\_deps

*Infer dependencies*


---

## Description

Infers the R packages that your R package depends on.

**Usage**

```
infer_deps(
  path = here::here(),
  which = c("Imports", "Suggests"),
  imports_thresh = 2,
  imports = NULL,
  suggests = c("testthat", "rmarkdown", "markdown", "knitr", "remotes", "knitr", "covr"),
  errors = c("reported", "fatal", "ignored"),
  dev = FALSE,
  progress = TRUE,
  add_newlines = FALSE
)
```

**Arguments**

path	The path to a (possibly multi-mode) R file, or a directory containing such files. By default, all files within the current working directory are checked, recursively.
which	Which types of dependencies to return.
imports_thresh	The minimum number of times that a package has to be called within your package to assign it as an Import. If is called less times than this threshold, it will instead be assigned as a Suggest, which means it will not be installed by default.
imports	R packages that are exempt from the <code>suggests_thresh</code> rule and are instead automatically assigned as Imports.
suggests	R packages that are exempt from the <code>suggests_thresh</code> rule and are instead automatically assigned as Suggests.
errors	How should errors that occur during dependency enumeration be handled? See <b>Errors</b> for more details.
dev	Boolean; include 'development' dependencies as well? That is, packages which may be required during development but are unlikely to be required during runtime for your project. By default, only runtime dependencies are returned.
progress	Boolean; report progress output while enumerating dependencies?
add_newlines	Prefix each package name with a newline character and two spaces. This is useful for formatting <i>DESCRIPTION</i> files.

**Value**

A character vector of R package names.

**Examples**

```
deps <- infer_deps()
```

---

is_gha	<i>Is GitHub Action</i>
--------	-------------------------

---

**Description**

Tests whether a function is currently being run within a GitHub Actions workflow or not.

**Usage**

```
is_gha(var = "GITHUB_ACTION", verbose = TRUE)
```

**Arguments**

var	Environmental variable to check.
verbose	Print messages.

**Source**

[GitHub Actions docs](#)

**Examples**

```
is_gha()
```

---

use_badges	<i>Use badges</i>
------------	-------------------

---

**Description**

Create one or more badges showing the status of your R package. Uses the package **badger**.

**Usage**

```
use_badges(  
  ref = NULL,  
  add_hex = TRUE,  
  add_actions = "rworkflows",  
  add_doi = NULL,  
  add_lifecycle = FALSE,  
  add_github_version = TRUE,  
  add_commit = TRUE,  
  add_code_size = TRUE,  
  add_license = TRUE,  
  add_authors = TRUE,  
  add_codecov = TRUE,  
  add_codecov_graphs = "icicle",  
  add_bioc_release = FALSE,  
  add_bioc_download_month = FALSE,  
  add_bioc_download_total = FALSE,  
  add_bioc_download_rank = FALSE,  
)
```

```

    add_cran_release = FALSE,
    add_cran_checks = FALSE,
    add_cran_download_month = FALSE,
    add_cran_download_total = FALSE,
    branch = "master",
    as_list = FALSE,
    sep = "\n",
    hex_height = 300,
    codecov_graph_width = 200,
    colors = list(github = "black", bioc = "green", cran = "black", default = "blue",
                  lifecycle = NULL),
    verbose = TRUE
  )

```

## Arguments

ref	Reference for a GitHub repository. If NULL (the default), the reference is determined by the URL field in the DESCRIPTION file.
add_hex	Add a hex sticker. If add_hex=TRUE, will assume the sticker is located at the following relative path: "inst/hex/hex.png". If add_hex is a character string, this will instead be used as the relative hex path (e.g. "/images/mysticker.png").
add_actions	The name of one or more GitHub Actions to show the status for with <a href="#">badge_github_actions</a> (e.g. c("rworkflows", "rworkflows_static")).
add_doi	Add the DOI of a given package or publication associated with the package using <a href="#">badge_doi</a> . Must be provided as a character string, e.g.: "10.1111/2041-210X.12628"
add_lifecycle	Add package lifecycle stage. If not FALSE, must be a character string indicating one of the following valid lifecycle stage: <ul style="list-style-type: none"> <li>• "stable"</li> <li>• "deprecated"</li> <li>• "superseded"</li> <li>• "experimental"</li> </ul> See <a href="https://lifecycle.r-lib.org">lifecycle.r-lib.org</a> for further details.
add_github_version	Add package version with <a href="#">badge_github_version</a> .
add_commit	Add the last GitHub repo commit date with <a href="#">badge_last_commit</a> .
add_code_size	Add code size with <a href="#">badge_code_size</a> .
add_license	Add license info with <a href="#">badge_license</a> .
add_authors	Add author names inferred from the DESCRIPTION file.
add_codecov	Add Codecov status with <a href="#">badge_codecov</a> . See the <a href="#">Codecov site</a> for more information about these badges.
add_codecov_graphs	Add Codecov graphs visualising results of code coverage tests. Options include: <ul style="list-style-type: none"> <li>• "sunburst"</li> <li>• "tree"</li> <li>• "icicle"</li> </ul> See the <a href="#">Codecov site</a> for more information about each plot type.
add_bioc_release	Add Bioc release version with <a href="#">badge_bioc_release</a> .

add_bioc_download_month	Add the number of Bioc downloads last month <a href="#">badge_bioc_download</a> .
add_bioc_download_total	Add the number of Bioc downloads total <a href="#">badge_bioc_download</a> .
add_bioc_download_rank	Add the download rank of the package on Bioc <a href="#">badge_bioc_download_rank</a> .
add_cran_release	Add Bioc release version with <a href="#">badge_cran_release</a> .
add_cran_checks	Add whether package is passing all checks on CRAN with <a href="#">badge_cran_checks</a> .
add_cran_download_month	Add the number of CRAN downloads last month <a href="#">badge_cran_download</a> .
add_cran_download_total	Add the number of CRAN downloads total <a href="#">badge_cran_download</a> .
branch	Name of the GitHub repository branch to use.
as_list	Return the header as a named list (TRUE), or a collapsed text string (default: FALSE).
sep	Character to separate each item in the list with using <a href="#">paste</a> .
hex_height	Height of the hex sticker in pixels (when add_hex=TRUE).
codecov_graph_width	Width of each Codecov graph in pixels (when add_codecov_graph!=FALSE).
colors	Colors to assign to each group of badges (when possible).
verbose	Print messages.

### Value

A named list of selected badges in markdown format.

### Examples

```
badges <- rworkflows::use_badges(ref = "neurogenomics/rworkflows")
```

---

use_dockerfile	<i>Use Dockerfile</i>
----------------	-----------------------

---

### Description

Creates a Docker file to be used with the GitHub Actions (GHA) workflows distributed by **rworkflows**.

### Usage

```
use_dockerfile(
  save_dir = here::here(),
  path = file.path(save_dir, "Dockerfile"),
  force_new = FALSE,
  show = FALSE,
  verbose = TRUE
)
```

**Arguments**

save_dir	Directory to save the Docker file to.
path	Path to the Docker file.
force_new	If a Docker file already exists, overwrite it (default: FALSE).
show	Print the contents of the Docker file in the R console.
verbose	Print messages.

**Value**

Path to Docker file.

**Examples**

```
path <- use_dockerfile(save_dir=tempdir())
```

---

use_issue_template	<i>Use Issue Template</i>
--------------------	---------------------------

---

**Description**

Creates one or more Issue Templates to be used in a GitHub repository.

**Usage**

```
use_issue_template(
  templates = c("bug_report.md", "feature_request.md"),
  save_dir = here::here(".github", "ISSUE_TEMPLATE"),
  path = file.path(save_dir, templates),
  force_new = FALSE,
  show = FALSE,
  verbose = TRUE
)
```

**Arguments**

templates	The names of templates to be used.
save_dir	Directory to save the Docker file to.
path	Path to the Docker file.
force_new	If a Docker file already exists, overwrite it (default: FALSE).
show	Print the contents of the Docker file in the R console.
verbose	Print messages.

**Value**

Path to Issue Templates.

**Examples**

```
path <- use_issue_template(save_dir=tempdir())
```

---

use_readme	<i>Use README</i>
------------	-------------------

---

### Description

Creates an rmarkdown README file that autofills using metadata from the R package *DESCRIPTION* file.

### Usage

```
use_readme(
  save_dir = here::here(),
  path = file.path(save_dir, "README.Rmd"),
  force_new = FALSE,
  show = FALSE,
  verbose = TRUE
)
```

### Arguments

save_dir	Directory to save the vignette file to.
path	Path to the vignette file.
force_new	If the file already exists, overwrite it (default: FALSE).
show	Print the contents of the vignette file in the R console.
verbose	Print messages.

### Value

Path to README file.

### Examples

```
## use default save_dir in practice
path <- use_readme(save_dir = tempdir())
```

---

use_vignette_docker	<i>Use vignette: Docker</i>
---------------------	-----------------------------

---

### Description

Creates a vignette rmarkdown file demonstrates how to create a Docker/Singularity image from a container stored in [Dockerhub](#).



## Usage

```
use_vignette_docker(  
  docker_org,  
  title = "Docker/Singularity Containers",  
  vignette_index_entry = "docker",  
  save_dir = here::here(),  
  path = file.path(save_dir, "vignettes", "docker.Rmd"),  
  output = "BiocStyle::html_document",  
  port_in = 8787,  
  port_out = 8900,  
  force_new = FALSE,  
  show = FALSE,  
  verbose = TRUE  
)
```

## Arguments

<code>docker_org</code>	DockerHub organization name. Can simply be your Dockerhub username instead.
<code>title</code>	Title of vignette.
<code>vignette_index_entry</code>	Index entry of the vignette, which is used when creating the navigation bar in the <b>pkgdown</b> site.
<code>save_dir</code>	Directory to save the vignette file to.
<code>path</code>	Path to the vignette file.
<code>output</code>	Vignette output style. Defaults to <a href="#">html_document</a> .
<code>port_in</code>	Port number to route into the docker container. See the <a href="#">Docker docs</a> for further details.
<code>port_out</code>	Port number to route out of docker container. See the <a href="#">Docker docs</a> for further details.
<code>force_new</code>	If the file already exists, overwrite it (default: FALSE).
<code>show</code>	Print the contents of the vignette file in the R console.
<code>verbose</code>	Print messages.

## Value

Path to vignette file.

## Examples

```
path <- use_vignette_docker(docker_org = "neurogenomicslab",  
                           ## use default save_dir in practice  
                           save_dir = tempdir())
```

---

`use_vignette_getstarted`*Use vignette: Get started*

---

## Description

Creates a "Get started" rmarkdown vignette file.

## Usage

```
use_vignette_getstarted(  
  package,  
  title = "Get started",  
  vignette_index_entry = package,  
  save_dir = here::here(),  
  path = file.path(save_dir, "vignettes", paste0(package, ".Rmd")),  
  output = "BiocStyle::html_document",  
  force_new = FALSE,  
  show = FALSE,  
  verbose = TRUE  
)
```

## Arguments

<code>package</code>	R package name.
<code>title</code>	Title of vignette.
<code>vignette_index_entry</code>	Index entry of the vignette, which is used when creating the navigation bar in the <b>pkgdown</b> site.
<code>save_dir</code>	Directory to save the vignette file to.
<code>path</code>	Path to the vignette file.
<code>output</code>	Vignette output style. Defaults to <a href="#">html_document</a> .
<code>force_new</code>	If the file already exists, overwrite it (default: FALSE).
<code>show</code>	Print the contents of the vignette file in the R console.
<code>verbose</code>	Print messages.

## Value

Path to vignette file.

## Examples

```
path <- use_vignette_getstarted(package = "mypackage",  
                                ## use default save_dir in practice  
                                save_dir = tempdir())
```

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use_workflow	<i>Use GitHub Actions workflow</i>
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## Description

Create workflow that calls an **rworkflows** **GitHub Actions (GHA)**

## Usage

```
use_workflow(
  name = "rworkflows",
  tag = "@master",
  on = c("push", "pull_request"),
  branches = c("master", "main", "devel", "RELEASE_**"),
  runners = construct_runners(),
  run_bioccheck = FALSE,
  run_rcmdcheck = TRUE,
  as_cran = TRUE,
  run_vignettes = TRUE,
  has_testthat = TRUE,
  run_covr = TRUE,
  run_pkgdown = TRUE,
  has_runit = FALSE,
  has_latex = FALSE,
  run_docker = FALSE,
  github_token = "${ secrets.PAT_GITHUB }",
  docker_user = NULL,
  docker_org = docker_user,
  docker_token = "${ secrets.DOCKER_TOKEN }",
  cache_version = "cache-v1",
  enable_act = FALSE,
  save_dir = here::here(".github", "workflows"),
  return_path = TRUE,
  force_new = FALSE,
  preview = FALSE,
  verbose = TRUE
)
```

## Arguments

name	Workflow name. <ul style="list-style-type: none"> <li>"rworkflows" A short workflow script that calls the GitHub action from the GitHub Marketplace. The action is continually updated so users do not need to worry about maintaining it.</li> <li>"rworkflows_static" A longer workflow scripts that explicitly copies all steps from the <b>rworkflows</b> action into a static file. Users may need to update this file themselves over time, though this does allow for a fully customisable workflow.</li> </ul>
tag	Which version of the rworkflows action to use. Can be a branch name on the <b>GitHub repository</b> (e.g. "@master"), or a <b>Release Tag</b> (e.g. "@v1").

on	GitHub trigger conditions.
branches	GitHub trigger branches.
runners	Runner configurations for multiple Operating Systems (OS), including R versions, Bioc versions, and container sources. Can use the <a href="#">construct_runners</a> functions to assist in constructing customized runners configurations.
run_bioccheck	Run Bioconductor checks using <code>BiocCheck::BiocCheck()</code> . Must pass in order to continue workflow.
run_rcmdcheck	Run R CMD checks using <code>rcmdcheck::rcmdcheck()</code> . Must pass in order to continue workflow.
as_cran	When running R CMD checks, use the '-as-cran' flag to apply CRAN standards
run_vignettes	Build and check R package vignettes.
has_testthat	Run unit tests and report results.
run_covr	Run code coverage tests and publish results to codecov.
run_pkgdown	Knit the <i>README.Rmd</i> (if available), build documentation website, and deploy to <i>gh-pages</i> branch.
has_runit	Run R Unit tests.
has_latex	Install a suite of LaTeX dependencies used for rendering Sweave (.rnw) and other documentation files.
run_docker	Whether to build and push a Docker container to DockerHub.
github_token	Token for the repo. Can be passed in using secrets.PAT_GITHUB .
docker_user	DockerHub username.
docker_org	DockerHub organization name. Is the same as docker_user by default.
docker_token	DockerHub token.
cache_version	Name of the cache subdirectory to be used when reinstalling software in GHA.
enable_act	Whether to add extra lines to the yaml to enable local workflow checking with <code>act</code> .
save_dir	Directory to save workflow to.
return_path	Return the path to the saved <i>yaml</i> workflow file (default: TRUE), or return the <i>yaml</i> object directly.
force_new	If the GHA workflow yaml already exists, overwrite with new one (default: FALSE).
preview	Print the yaml file to the R console.
verbose	Print messages.

**Value**

Path or yaml object.

**Source**

Issue reading in "on:"/"y","n" elements.

Issue writing "on:" as "'as':"

**Examples**

```
path <- use_workflow(save_dir = file.path(tempdir(), ".github", "workflows"))
```

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