

# Package ‘systemPipeRdata’

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**Title** systemPipeRdata: Workflow templates and sample data

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**biocViews** Genetics, Infrastructure, DataImport, Sequencing, RNASeq, ChIPSeq, MethylSeq, SNP, GeneExpression, Coverage, GeneSetEnrichment, Alignment, QualityControl, ImmunoOncology, RiboSeq, WorkflowStep

**Description** systemPipeRdata complements the systemPipeR workflow management system (WMS) by offering a collection of pre-designed data analysis workflow templates. These templates are easily accessible and can be readily loaded onto a user's system with a single command. Once loaded, the WMS can immediately utilize these templates for efficient end-to-end analysis, serving a wide range of data analysis needs.

**Depends** R (>= 3.6.0)

**Imports** methods, Biostrings, BiocGenerics, jsonlite, remotes

**Suggests** GenomicFeatures, GenomicRanges, IRanges, Rsamtools, ShortRead, rtracklayer, RUnit, BiocStyle, knitr, rmarkdown, systemPipeR, kableExtra, magrittr, dplyr

**VignetteBuilder** knitr

**License** Artistic-2.0

**NeedsCompilation** no

**URL** <https://github.com/tgirke/systemPipeRdata>, <https://systempipe.org/>

**git\_url** <https://git.bioconductor.org/packages/systemPipeRdata>

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availableWF	<i>List Available Workflows Templates at systemPipeRdata package</i>
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### Description

This function checks the workflow templates availability from systemPipeRdata package and also from [systemPipeR Organization](#) on GitHub.

### Usage

```
availableWF(github = FALSE)
```

### Arguments

github	logical. If TRUE, it will return current experimental workflow templates available on systemPipeR Organization.
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### Details

Internally, this function uses the GitHub API, and there is an access limit per hour. For more details, please check: `system("curl -i https://api.github.com/users/<username>")`.

### Value

Workflow templates are printed on the console, and also return an invisible list with the names of the workflows templates available at systemPipeRdata package. If `github = TRUE`, it will return an additional data.frame with current workflow templates available on systemPipeR Organization in the list.

### Note

We are assuming that workflow templates repositories under [systemPipeR Organization](#) content the keyword "Workflow Template" on the Description section and "Topics" section, we expected "systempiper" and "release" or "development" words.

### Author(s)

Daniela Cassol

### See Also

[genWorkenvir](#).

**Examples**

```

availableWF()
## Not run:
## List Workflow Templates from \code{systemPipeR} Organization
availableWF(github = TRUE)

## End(Not run)

```

---

genWorkenvir

*Generate workflow templates*


---

**Description**

Generates workflow templates for `systemPipeR` package. The template environments contain a predefined directory structure along with run parameter files and sample data. The structure of the workflow templates and the sample data are described in all details in the Overview Vignette of the [systemPipeR package](#).

**Usage**

```
genWorkenvir(workflow, mydirname=NULL, bam=FALSE)
```

**Arguments**

workflow	character string of workflow templates to be generated. Supported values can be checked with the <code>\link{availableWF}()</code> function.
mydirname	Specifies the name of the workflow directory. The default NULL uses the name of the chosen workflow. An error is issued if a directory of the same name and path exists already.
bam	If <code>bam=TRUE</code> pregenerated short read alignment (BAM) files will be included in the <code>results</code> directory of the workflow environment. Note, these BAM files have been generated with the HISAT2 aligner using the FASTQ files provided in the <code>data</code> directory. The default <code>bam=FALSE</code> omits this step meaning no BAM files will be copied into the <code>results</code> directory.

**Details**

Check the output of `\link{availableWF}()` to the current workflow templates available on `systemPipeR` Organization.

For an `interactive()` session, the `readline()` function provides the option choose between proceeding or not, through options: `yes` or `no`. For non-interactive use, if there is no package install, the option `yes` will be selected.

**Value**

Workflow directory containing sample data and parameter files along with the following subdirectories:

param/	stores parameter files
data/	stores input data

results/            stores output results

For more details, please consult the Overview Vignette (HTML) of the systemPipeR package (<http://bioconductor.org/packages/systemPipeR>).

### Author(s)

Thomas Girke and Daniela Cassol

### Examples

```
## Return location of sample data
samplepaths <- pathList()
## Not run:
## Generate varseq workflow environment
genWorkenvir(workflow="varseq", mydirname=NULL)
setwd("varseq")

## End(Not run)
```

---

getSubsetReads            *Subsetting fastq data*

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### Description

Returns subsets of fastq files data based on specific mapping regions or list of genes or GRanges object.

### Usage

```
getSubsetReads(args,
  geneList = NULL,
  gr = NULL,
  MappingRegion = 1:1e+05,
  sample_range = 90000:1e+05,
  truncate_refs = TRUE,
  id_read_number = TRUE,
  annotation = "data/tair10.gff",
  reference = "data/tair10.fasta",
  annot_outname = "tair10_sub.gff",
  ref_outname = "tair10_sub.fasta",
  outdir = "data/subset/",
  silent = FALSE
)
```

### Arguments

args	object of class SYSargs2.
geneList	selected genes list to retrieve the reads from the fastq file.
gr	an object containing genomic ranges to retrieve the reads from the fastq file.
MappingRegion	integers ranges of start and end of chromosome position to retrieve the reads from the fastq file.

sample_range	random range to subsetted the fastq file.
truncate_refs	logical. If TRUE it will generate reference genome and annotation subset file.
id_read_number	if fastq file contains sequence name with read number ( <code>\$ri --define-seq '@\$sn[_\$rn]/\$ri '</code> ).
annotation	path to annotation file.
reference	path to reference genome.
annot_outname	character name of the annotation output file.
ref_outname	character name of the reference genome output file.
outdir	path to output directory.
silent	if set to TRUE, all messages returned by the function will be suppressed.

**Value**

Workflow directory containing sample data and parameter files along with the following subdirectories:

param/	stores parameter files
data/	stores input data
results/	stores output results

For more details, please consult the Overview Vignette (HTML) of the systemPipeR package (<http://bioconductor.org/packages/systemPipeR>).

**Author(s)**

Thomas Girke, Shiyuan Guo and Daniela Cassol

**Examples**

```
## Not run:
getSubsetReads(args, MappingRegion = 1:900, sample_range = 800:900, outdir = "data/subset/", silent = FALSE)
getSubsetReads(args, MappingRegion = 1:900, sample_range = NULL, outdir = "data/subset/", silent = FALSE)

## End(Not run)
```

---

pathList	<i>Return location of sample data</i>
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**Description**

Function to return paths to sample data provided by sytemPipeRdata package.

**Usage**

```
pathList()
```

**Value**

list

**Author(s)**

Thomas Girke

**Examples**

```
samplepaths <- pathList()
```

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