

# Package ‘mlr3resampling’

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**Type** Package

**Title** Resampling Algorithms for 'mlr3' Framework

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**Description** A supervised learning algorithm inputs a train set, and outputs a prediction function, which can be used on a test set. If each data point belongs to a subset (such as geographic region, year, etc), then how do we know if subsets are similar enough so that we can get accurate predictions on one subset, after training on Other subsets? And how do we know if training on All subsets would improve prediction accuracy, relative to training on the Same subset? SOAK, Same/Other/All K-fold cross-validation, <[doi:10.48550/arXiv.2410.08643](https://doi.org/10.48550/arXiv.2410.08643)> can be used to answer these questions, by fixing a test subset, training models on Same/Other/All subsets, and then comparing test error rates (Same versus Other and Same versus All). Also provides code for estimating how many train samples are required to get accurate predictions on a test set.

**License** LGPL-3

**URL** <https://github.com/tdhock/mlr3resampling>

**BugReports** <https://github.com/tdhock/mlr3resampling/issues>

**Imports** data.table, R6, checkmate, paradox, mlr3 (>= 0.21.1), mlr3misc, batchtools, filelock

**Suggests** ggplot2, animint2, mlr3tuning, lgr, future, testthat, knitr, markdown, nc, rpart, directlabels, mlr3torch, torch

**VignetteBuilder** knitr

**NeedsCompilation** no

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AZtrees

*Arizona Trees*

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**Description**

Classification data set with polygons (groups which should not be split in CV) and subsets (region3 or region4).

**Usage**

```
data("AZtrees")
```

**Format**

A data frame with 5956 observations on the following 25 variables.

region3 a character vector  
region4 a character vector  
polygon a numeric vector  
y a character vector  
ycoord latitude  
xcoord longitude  
SAMPLE\_1 a numeric vector  
SAMPLE\_2 a numeric vector  
SAMPLE\_3 a numeric vector  
SAMPLE\_4 a numeric vector  
SAMPLE\_5 a numeric vector  
SAMPLE\_6 a numeric vector  
SAMPLE\_7 a numeric vector  
SAMPLE\_8 a numeric vector  
SAMPLE\_9 a numeric vector  
SAMPLE\_10 a numeric vector  
SAMPLE\_11 a numeric vector  
SAMPLE\_12 a numeric vector  
SAMPLE\_13 a numeric vector  
SAMPLE\_14 a numeric vector  
SAMPLE\_15 a numeric vector  
SAMPLE\_16 a numeric vector  
SAMPLE\_17 a numeric vector  
SAMPLE\_18 a numeric vector  
SAMPLE\_19 a numeric vector  
SAMPLE\_20 a numeric vector  
SAMPLE\_21 a numeric vector

**Source**

Paul Nelson Arellano, paul.arellano@nau.edu

**Examples**

```
data(AZtrees)
task.obj <- mlr3::TaskClassif$new("AZtrees3", AZtrees, target="y")
task.obj$col_roles$feature <- grep("SAMPLE", names(AZtrees), value=TRUE)
task.obj$col_roles$group <- "polygon"
task.obj$col_roles$subset <- "region3"
str(task.obj)
same_other_sizes_cv <- mlr3resampling::ResamplingSameOtherSizesCV$new()
same_other_sizes_cv$instantiate(task.obj)
same_other_sizes_cv$instance$iteration.dt
```

---

proj\_compute

---

*Compute resampling results in a project*


---

**Description**

proj\_compute() looks in grid\_jobs.csv for a row with status=="not started", and picks the first one to work on, updating status="started". After having run train() and predict(), a data table with one row is saved to an RDS file in the grid\_jobs directory, and status="done" is updated. proj\_compute\_until\_done() keeps doing that in a while loop.

**Usage**

```
proj_compute(proj_dir, verbose = FALSE)
proj_compute_until_done(proj_dir, verbose = FALSE)
```

**Arguments**

proj_dir	Project directory created by <a href="#">proj_grid</a> .
verbose	Logical: print messages?

**Details**

If everything goes well, the user should not need to run this function. Instead, the user runs [proj\\_submit](#) as Step 2 out of the typical 3 step pipeline (init grid, submit, read results). proj\_compute can sometimes be useful for testing or debugging the submit step, since it runs one split at a time.

**Value**

Data table with one row of results.

**Author(s)**

Toby Dylan Hocking

## Examples

```

N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=factor(rep(c("Alice","Bob"), each=0.5*N))
)
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^as.integer(person))
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  task.dt <- data.table(reg.dt)[
    , y := f(x,person)+rnorm(N, sd=0.5)
  ][]
  task.obj <- mlr3::TaskRegr$new(
    pattern, task.dt, target="y")
  task.obj$col_roles$feature <- "x"
  task.obj$col_roles$stratum <- "person"
  task.obj$col_roles$subset <- "person"
  reg.task.list[[pattern]] <- task.obj
}
reg.learner.list <- list(
  featureless=mlr3::LearnerRegrFeatureless$new())
if(requireNamespace("rpart")){
  reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
}

pkg.proj.dir <- tempfile()
mlr3resampling::proj_grid(
  pkg.proj.dir,
  reg.task.list,
  reg.learner.list,
  SOAK,
  order_jobs = function(DT)1:2, # for CRAN.
  score_args=mlr3::msrs(c("regr.rmse", "regr.mae")))
mlr3resampling::proj_compute(pkg.proj.dir)
fread(file.path(pkg.proj.dir, "grid_jobs.csv"))
mlr3resampling::proj_compute_until_done(pkg.proj.dir)
fread(file.path(pkg.proj.dir, "grid_jobs.csv"))

```

---

proj\_grid

*Initialize a new project grid table*

---

## Description

A project grid consists of all combinations of tasks, learners, resampling types, and resampling iterations, to be computed in parallel. This function creates a project directory with files to describe

the grid.

## Usage

```
proj_grid(
  proj_dir, tasks, learners, resamplings,
  order_jobs = NULL, score_args = NULL,
  save_learner = FALSE, save_pred = FALSE)
```

## Arguments

proj_dir	Path to directory to create.
tasks	List of Tasks, or a single Task.
learners	List of Learners, or a single Learner.
resamplings	List of Resamplings, or a single Resampling.
order_jobs	Function which takes split table as input, and returns integer vector of row numbers of the split table to write to <code>grid_jobs.csv</code> , which is how worker processes determine what work to do next (smaller numbers have higher priority). Default NULL means to keep default order.
score_args	Passed to <code>pred\$score()</code> .
save_learner	Function to process Learner, after training/prediction, but before saving result to disk. For interpreting complex models, you should write a function that returns only the parts of the model that you need (and discards the other parts which would take up disk space for no reason). Default FALSE means to not keep it (always returns NULL). TRUE means to keep it without any special processing.
save_pred	Function to process Prediction before saving to disk. Default FALSE means to not keep it (always returns NULL). TRUE means to keep it without any special processing.

## Details

This is Step 1 out of the typical 3 step pipeline (init grid, submit, read results). It creates a `grid_jobs.csv` table which has a column `status`; each row is initialized to "not started" or "done", depending on whether the corresponding result RDS file exists already.

## Value

Data table of splits to be processed (same as table saved to `grid_jobs.rds`).

## Author(s)

Toby Dylan Hocking

## Examples

```

N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=factor(rep(c("Alice", "Bob"), each=0.5*N))
)
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^as.integer(person))
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  task.dt <- data.table(reg.dt)[
    , y := f(x,person)+rnorm(N, sd=0.5)
  ][]
  task.obj <- mlr3::TaskRegr$new(
    pattern, task.dt, target="y")
  task.obj$col_roles$feature <- "x"
  task.obj$col_roles$stratum <- "person"
  task.obj$col_roles$subset <- "person"
  reg.task.list[[pattern]] <- task.obj
}
reg.learner.list <- list(
  featureless=mlr3::LearnerRegrFeatureless$new())
if(requireNamespace("rpart")){
  reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
}

pkg.proj.dir <- tempfile()
mlr3resampling::proj_grid(
  pkg.proj.dir,
  reg.task.list,
  reg.learner.list,
  SOAK,
  score_args=mlr3::msrs(c("regr.rmse", "regr.mae")))
mlr3resampling::proj_compute(pkg.proj.dir)
fread(file.path(pkg.proj.dir, "grid_jobs.csv"))

```

---

proj\_results

Combine and save results in a project

---

## Description

proj\_results globs the RDS result files in the project directory, and combines them into a result table via rbindlist(). proj\_results\_save saves that result table to results.rds and results.csv.

**Usage**

```
proj_results(proj_dir)
proj_results_save(proj_dir)
```

**Arguments**

proj\_dir            Project directory created via [proj\\_grid](#).

**Details**

This is Step 3 out of the typical 3 step pipeline (init grid, submit, read results). Actually, if step 2 worked as intended, the last [proj\\_compute](#) calls [proj\\_results\\_save](#), which saves up to three result files to disk that you can read directly:

results.csv contains test measures for each split, and can be read via `fread()`

results.rds contains additional list columns for learner and pred (useful for model interpretation), and can be read via `readRDS()`

learners.csv only exists if learner column is a data frame, in which case it contains the atomic columns, along with meta-data describing each split.

**Value**

`proj_results` returns a data table with all columns, whereas `proj_results_save` returns the same table with only atomic columns.

**Author(s)**

Toby Dylan Hocking

**Examples**

```
N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=factor(rep(c("Alice", "Bob"), each=0.5*N)))
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^as.integer(person))
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  task.dt <- data.table(reg.dt[,
    , y := f(x,person)+rnorm(N, sd=0.5)
  ])
  task.obj <- mlr3::TaskRegr$new(
    pattern, task.dt, target="y")
  task.obj$col_roles$feature <- "x"
```



```

    task.obj$col_roles$stratum <- "person"
    task.obj$col_roles$subset <- "person"
    reg.task.list[[pattern]] <- task.obj
  }
  reg.learner.list <- list(
    featureless=mlr3::LearnerRegrFeatureless$new()
  )
  if(requireNamespace("rpart")){
    reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
  }

  pkg.proj.dir <- tempfile()
  mlr3resampling::proj_grid(
    pkg.proj.dir,
    reg.task.list,
    reg.learner.list,
    SOAK,
    order_jobs = function(DT)1:2, # for CRAN.
    score_args=mlr3::msrs(c("regr.rmse", "regr.mae")))
  mlr3resampling::proj_compute_until_done(pkg.proj.dir)
  fread(file.path(pkg.proj.dir, "results.csv"))

```

proj\_submit

*Submit resampling split jobs in parallel*

## Description

Before running this function, you should define `cluster.functions` in your `~/.batchtools.conf.R` file. It makes a batchtools registry, then runs `batchtools::batchMap()` and `batchtools::submitJobs()`; each iteration runs [proj\\_compute\\_until\\_done](#).

## Usage

```

proj_submit(
  proj_dir, tasks = 2, hours = 1, gigabytes = 1,
  verbose = FALSE, cluster.functions = NULL)

```

## Arguments

<code>proj_dir</code>	Project directory created via <a href="#">proj_grid</a> .
<code>tasks</code>	Positive integer: number of batchtools jobs, translated into one SLURM job array with this number of tasks.
<code>hours</code>	Hours of walltime to ask the SLURM scheduler.
<code>gigabytes</code>	Gigabytes of memory to ask the SLURM scheduler.
<code>verbose</code>	Logical: print messages?
<code>cluster.functions</code>	Cluster functions from batchtools, useful for testing.

**Details**

This is Step 2 out of the typical 3 step pipeline (init grid, submit, read results).

**Value**

The batchtools registry.

**Author(s)**

Toby Dylan Hocking

**Examples**

```

N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=factor(rep(c("Alice","Bob"), each=0.5*N)))
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^as.integer(person))
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  task.dt <- data.table(reg.dt)[
    , y := f(x,person)+rnorm(N, sd=0.5)
  ][]
  task.obj <- mlr3::TaskRegr$new(
    pattern, task.dt, target="y")
  task.obj$col_roles$feature <- "x"
  task.obj$col_roles$stratum <- "person"
  task.obj$col_roles$subset <- "person"
  reg.task.list[[pattern]] <- task.obj
}
reg.learner.list <- list(
  featureless=mlr3::LearnerRegrFeatureless$new())
if(requireNamespace("rpart")){
  reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
}

pkg.proj.dir <- tempfile()
mlr3resampling::proj_grid(
  pkg.proj.dir,
  reg.task.list,
  reg.learner.list,
  SOAK,
  order_jobs = function(DT)1:2, # for CRAN.
  score_args=mlr3::msrs(c("regr.rmse", "regr.mae")))
mlr3resampling::proj_submit(pkg.proj.dir)
batchtools::waitForJobs()

```

```
fread(file.path(pkg.proj.dir, "results.csv"))
```

---

pvalue	<i>P-values for comparing Same/Other/All training</i>
--------	---

---

## Description

Same/Other/All K-fold cross-validation (SOAK) results in K measures of test error/accuracy. This function computes P-values (two-sided T-test) between Same and All/Other.

## Usage

```
pvalue(score_in, value.var = NULL, digits=3)
```

## Arguments

score_in	Data table output from <a href="#">score</a> .
value.var	Name of column to use as the evaluation metric in T-test. Default NULL means to use the first column matching "classif regr".
digits	Number of decimal places to show for mean and standard deviation.

## Value

List of class "pvalue" with named elements value.var, stats, pvalues.

## Author(s)

Toby Dylan Hocking

## Examples

```
N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=rep(1:2, each=0.5*N))
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^person)
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  yname <- paste0("y_",pattern)
  reg.dt[, (yname) := f(x,person)+rnorm(N, sd=0.5)][]
  task.dt <- reg.dt[, c("x","person",yname), with=FALSE]
  task.obj <- mlr3::TaskRegr$new(
```

```

    pattern, task.dt, target=yname)
  task.obj$col_roles$stratum <- "person"
  task.obj$col_roles$subset <- "person"
  reg.task.list[[pattern]] <- task.obj
}
reg.learner.list <- list(
  mlr3::LearnerRegrFeatureless$new()
)
if(requireNamespace("rpart")){
  reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
}
(bench.grid <- mlr3::benchmark_grid(
  reg.task.list,
  reg.learner.list,
  SOAK))
bench.result <- mlr3::benchmark(bench.grid)
bench.score <- mlr3resampling::score(bench.result, mlr3::msr("regr.rmse"))
bench.plist <- mlr3resampling::pvalue(bench.score)
plot(bench.plist)

```

---

ResamplingSameOtherCV *Resampling for comparing training on same or other subsets*

---

## Description

[ResamplingSameOtherCV](#) defines how a task is partitioned for resampling, for example in [resample\(\)](#) or [benchmark\(\)](#).

Resampling objects can be instantiated on a [Task](#), which should define at least one subset variable.

After instantiation, sets can be accessed via `$train_set(i)` and `$test_set(i)`, respectively.

## Details

This provides an implementation of SOAK, Same/Other/All K-fold cross-validation. After instantiation, this class provides information in `$instance` that can be used for visualizing the splits, as shown in the vignette. Most typical machine learning users should instead use [ResamplingSameOtherSizesCV](#), which does not support these visualization features, but provides other relevant machine learning features, such as group role, which is not supported by [ResamplingSameOtherCV](#).

A supervised learning algorithm inputs a train set, and outputs a prediction function, which can be used on a test set. If each data point belongs to a subset (such as geographic region, year, etc), then how do we know if it is possible to train on one subset, and predict accurately on another subset? Cross-validation can be used to determine the extent to which this is possible, by first assigning fold IDs from 1 to K to all data (possibly using stratification, usually by subset and label). Then we loop over test sets (subset/fold combinations), train sets (same subset, other subsets, all subsets), and compute test/prediction accuracy for each combination. Comparing test/prediction accuracy between same and other, we can determine the extent to which it is possible (perfect if same/other have similar test accuracy for each subset; other is usually somewhat less accurate than same; other can be just as bad as featureless baseline when the subsets have different patterns).

## Stratification

[ResamplingSameOtherCV](#) supports stratified sampling. The stratification variables are assumed to be discrete, and must be stored in the [Task](#) with column role "stratum". In case of multiple stratification variables, each combination of the values of the stratification variables forms a stratum.

## Grouping

[ResamplingSameOtherCV](#) does not support grouping of observations that should not be split in cross-validation. See [ResamplingSameOtherSizesCV](#) for another sampler which does support both group and subset roles.

## Subsets

The subset variable is assumed to be discrete, and must be stored in the [Task](#) with column role "subset". The number of cross-validation folds K should be defined as the fold parameter. In each subset, there will be about an equal number of observations assigned to each of the K folds. The assignments are stored in `$instance$id.dt`. The train/test splits are defined by all possible combinations of test subset, test fold, and train subsets (Same/Other/All). The splits are stored in `$instance$iteration.dt`.

## Methods

### Public methods:

- [Resampling\\$new\(\)](#)
- [Resampling\\$train\\_set\(\)](#)
- [Resampling\\$test\\_set\(\)](#)

**Method** `new()`: Creates a new instance of this [R6](#) class.

*Usage:*

```
Resampling$new(
  id,
  param_set = ps(),
  duplicated_ids = FALSE,
  label = NA_character_,
  man = NA_character_
)
```

*Arguments:*

```
id (character(1))
  Identifier for the new instance.
param_set (paradox::ParamSet)
  Set of hyperparameters.
duplicated_ids (logical(1))
  Set to TRUE if this resampling strategy may have duplicated row ids in a single training set
  or test set.
label (character(1))
  Label for the new instance.
```

`man` (character(1))

String in the format `[pkg]::[topic]` pointing to a manual page for this object. The referenced help package can be opened via method `$help()`.

**Method** `train_set()`: Returns the row ids of the *i*-th training set.

*Usage:*

`Resampling$train_set(i)`

*Arguments:*

*i* (integer(1))

Iteration.

*Returns:* (integer()) of row ids.

**Method** `test_set()`: Returns the row ids of the *i*-th test set.

*Usage:*

`Resampling$test_set(i)`

*Arguments:*

*i* (integer(1))

Iteration.

*Returns:* (integer()) of row ids.

### See Also

- arXiv paper <https://arxiv.org/abs/2410.08643> describing SOAK algorithm.
- Articles <https://github.com/tdhock/mlr3resampling/wiki/Articles>
- Package **mlr3** for standard [Resampling](#), which does not support comparing train on Same/Other/All subsets.
- `vignette(package="mlr3resampling")` for more detailed examples.

### Examples

```
same_other <- mlr3resampling::ResamplingSameOtherCV$new()
same_other$param_set$values$folds <- 5
```

---

ResamplingSameOtherSizesCV

*Resampling for comparing train subsets and sizes*

---

### Description

[ResamplingSameOtherSizesCV](#) defines how a task is partitioned for resampling, for example in [resample\(\)](#) or [benchmark\(\)](#).

Resampling objects can be instantiated on a [Task](#), which can use the subset role.

After instantiation, sets can be accessed via `$train_set(i)` and `$test_set(i)`, respectively.

## Details

This is an implementation of SOAK, Same/Other/All K-fold cross-validation. A supervised learning algorithm inputs a train set, and outputs a prediction function, which can be used on a test set. If each data point belongs to a subset (such as geographic region, year, etc), then how do we know if it is possible to train on one subset, and predict accurately on another subset? Cross-validation can be used to determine the extent to which this is possible, by first assigning fold IDs from 1 to K to all data (possibly using stratification, usually by subset and label). Then we loop over test sets (subset/fold combinations), train sets (same subset, other subsets, all subsets), and compute test/prediction accuracy for each combination. Comparing test/prediction accuracy between same and other, we can determine the extent to which it is possible (perfect if same/other have similar test accuracy for each subset; other is usually somewhat less accurate than same; other can be just as bad as featureless baseline when the subsets have different patterns).

This class has more parameters/potential applications than [ResamplingSameOtherCV](#) and [ResamplingVariableSizeTrainC](#) which are older and should only be preferred for visualization purposes.

## Stratification

[ResamplingSameOtherSizesCV](#) supports stratified sampling. The stratification variables are assumed to be discrete, and must be stored in the [Task](#) with column role "stratum". In case of multiple stratification variables, each combination of the values of the stratification variables forms a stratum.

## Grouping

[ResamplingSameOtherSizesCV](#) supports grouping of observations that will not be split in cross-validation. The grouping variable is assumed to be discrete, and must be stored in the [Task](#) with column role "group".

## Subsets

[ResamplingSameOtherSizesCV](#) supports training on different subsets of observations. The subset variable is assumed to be discrete, and must be stored in the [Task](#) with column role "subset".

## Parameters

The number of cross-validation folds K should be defined as the `fold` parameter, default 3.

The number of random seeds for down-sampling should be defined as the `seeds` parameter, default 1.

The ratio for down-sampling should be defined as the `ratio` parameter, default 0.5. The min size of same and other sets is repeatedly multiplied by this ratio, to obtain smaller sample sizes.

The number of down-sampling sizes/multiplications should be defined as the `sizes` parameter, which can also take two special values: default -1 means no down-sampling at all, and 0 means only down-sampling to the sizes of the same/other sets.

The `ignore_subset` parameter should be either TRUE or FALSE (default), whether to ignore the subset role. TRUE only creates splits for same subset (even if task defines subset role), and is useful for subtrain/validation splits (hyper-parameter learning). Note that this feature will work on a task with both stratum and group roles (unlike [ResamplingCV](#)).

The subsets parameter should specify the train subsets of interest: "S" (same), "O" (other), "A" (all), "SO", "SA", "SOA" (default).

In each subset, there will be about an equal number of observations assigned to each of the K folds. The train/test splits are defined by all possible combinations of test subset, test fold, train subsets (same/other/all), down-sampling sizes, and random seeds. The splits are stored in `$instance$iteration.dt`.

## Methods

### Public methods:

- `Resampling$new()`
- `Resampling$train_set()`
- `Resampling$test_set()`

**Method** `new()`: Creates a new instance of this R6 class.

*Usage:*

```
Resampling$new(
  id,
  param_set = ps(),
  duplicated_ids = FALSE,
  label = NA_character_,
  man = NA_character_
)
```

*Arguments:*

```
id (character(1))
  Identifier for the new instance.
param_set (paradox::ParamSet)
  Set of hyperparameters.
duplicated_ids (logical(1))
  Set to TRUE if this resampling strategy may have duplicated row ids in a single training set
  or test set.
label (character(1))
  Label for the new instance.
man (character(1))
  String in the format [pkg]::[topic] pointing to a manual page for this object. The refer-
  enced help package can be opened via method $help().
```

**Method** `train_set()`: Returns the row ids of the i-th training set.

*Usage:*

```
Resampling$train_set(i)
```

*Arguments:*

```
i (integer(1))
  Iteration.
```

*Returns:* `(integer())` of row ids.

**Method** `test_set()`: Returns the row ids of the i-th test set.



*Usage:*

```
Resampling$test_set(i)
```

*Arguments:*

*i* (integer(1))  
Iteration.

*Returns:* (integer()) of row ids.

## See Also

- arXiv paper <https://arxiv.org/abs/2410.08643> describing SOAK algorithm.
- Articles <https://github.com/tdhock/mlr3resampling/wiki/Articles>
- Package **mlr3** for standard [Resampling](#), which does not support comparing train on Same/Other/All subsets.
- `vignette(package="mlr3resampling")` for more detailed examples.

## Examples

```
same_other_sizes <- mlr3resampling::ResamplingSameOtherSizesCV$new()
same_other_sizes$param_set$values$folds <- 5
```

---

ResamplingVariableSizeTrainCV

*Resampling for comparing training on same or other groups*

---

## Description

[ResamplingVariableSizeTrainCV](#) defines how a task is partitioned for resampling, for example in [resample\(\)](#) or [benchmark\(\)](#).

Resampling objects can be instantiated on a [Task](#).

After instantiation, sets can be accessed via `$train_set(i)` and `$test_set(i)`, respectively.

## Details

A supervised learning algorithm inputs a train set, and outputs a prediction function, which can be used on a test set. How many train samples are required to get accurate predictions on a test set? Cross-validation can be used to answer this question, with variable size train sets.

## Stratification

[ResamplingVariableSizeTrainCV](#) supports stratified sampling. The stratification variables are assumed to be discrete, and must be stored in the [Task](#) with column role "stratum". In case of multiple stratification variables, each combination of the values of the stratification variables forms a stratum.

## Grouping

`ResamplingVariableSizeTrainCV` does not support grouping of observations.

## Hyper-parameters

The number of cross-validation folds should be defined as the `fold` parameter.

For each fold ID, the corresponding observations are considered the test set, and a variable number of other observations are considered the train set.

The `random_seeds` parameter controls the number of random orderings of the train set that are considered.

For each random order of the train set, the `min_train_data` parameter controls the size of the smallest stratum in the smallest train set considered.

To determine the other train set sizes, we use an equally spaced grid on the log scale, from `min_train_data` to the largest train set size (all data not in test set). The number of train set sizes in this grid is determined by the `train_sizes` parameter.

## Methods

### Public methods:

- `Resampling$new()`
- `Resampling$train_set()`
- `Resampling$test_set()`

**Method** `new()`: Creates a new instance of this R6 class.

*Usage:*

```
Resampling$new(
  id,
  param_set = ps(),
  duplicated_ids = FALSE,
  label = NA_character_,
  man = NA_character_
)
```

*Arguments:*

`id` (character(1))

Identifier for the new instance.

`param_set` (`paradox::ParamSet`)

Set of hyperparameters.

`duplicated_ids` (logical(1))

Set to TRUE if this resampling strategy may have duplicated row ids in a single training set or test set.

`label` (character(1))

Label for the new instance.

`man` (character(1))

String in the format `[pkg]::[topic]` pointing to a manual page for this object. The referenced help package can be opened via method `$help()`.

**Method** `train_set()`: Returns the row ids of the i-th training set.

*Usage:*

```
Resampling$train_set(i)
```

*Arguments:*

`i` (`integer(1)`)  
Iteration.

*Returns:* (`integer()`) of row ids.

**Method** `test_set()`: Returns the row ids of the i-th test set.

*Usage:*

```
Resampling$test_set(i)
```

*Arguments:*

`i` (`integer(1)`)  
Iteration.

*Returns:* (`integer()`) of row ids.

## Examples

```
(var_sizes <- mlr3resampling::ResamplingVariableSizeTrainCV$new())
```

---

score	<i>Score benchmark results</i>
-------	--------------------------------

---

## Description

Computes a data table of scores.

## Usage

```
score(bench.result, ...)
```

## Arguments

`bench.result`    Output of `benchmark()`.  
`...`            Additional arguments to pass to `bench.result$score`, for example measures.

## Value

data table with scores.

## Author(s)

Toby Dylan Hocking

## Examples

```

N <- 80
library(data.table)
set.seed(1)
reg.dt <- data.table(
  x=runif(N, -2, 2),
  person=rep(1:2, each=0.5*N))
reg.pattern.list <- list(
  easy=function(x, person)x^2,
  impossible=function(x, person)(x^2)*(-1)^person)
SOAK <- mlr3resampling::ResamplingSameOtherSizesCV$new()
reg.task.list <- list()
for(pattern in names(reg.pattern.list)){
  f <- reg.pattern.list[[pattern]]
  yname <- paste0("y_",pattern)
  reg.dt[, (yname) := f(x,person)+rnorm(N, sd=0.5)][]
  task.dt <- reg.dt[, c("x", "person", yname), with=FALSE]
  task.obj <- mlr3::TaskRegr$new(
    pattern, task.dt, target=yname)
  task.obj$col_roles$stratum <- "person"
  task.obj$col_roles$subset <- "person"
  reg.task.list[[pattern]] <- task.obj
}
reg.learner.list <- list(
  mlr3::LearnerRegrFeatureless$new())
if(requireNamespace("rpart")){
  reg.learner.list$rpart <- mlr3::LearnerRegrRpart$new()
}
(bench.grid <- mlr3::benchmark_grid(
  reg.task.list,
  reg.learner.list,
  SOAK))
bench.result <- mlr3::benchmark(bench.grid)
bench.score <- mlr3resampling::score(bench.result, mlr3::msr("regr.rmse"))
plot(bench.score)

```

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