

# Package ‘metasnf’

February 4, 2025

**Title** Meta Clustering with Similarity Network Fusion

**Version** 2.0.0

**Description** Framework to facilitate patient subtyping with similarity network fusion and meta clustering. The similarity network fusion (SNF) algorithm was introduced by Wang et al. (2014) in <[doi:10.1038/nmeth.2810](https://doi.org/10.1038/nmeth.2810)>. SNF is a data integration approach that can transform high-dimensional and diverse data types into a single similarity network suitable for clustering with minimal loss of information from each initial data source. The meta clustering approach was introduced by Caruana et al. (2006) in <[doi:10.1109/ICDM.2006.103](https://doi.org/10.1109/ICDM.2006.103)>. Meta clustering involves generating a wide range of cluster solutions by adjusting clustering hyperparameters, then clustering the solutions themselves into a manageable number of qualitatively similar solutions, and finally characterizing representative solutions to find ones that are best for the user's specific context. This package provides a framework to easily transform multimodal data into a wide range of similarity network fusion-derived cluster solutions as well as to visualize, characterize, and validate those solutions. Core package functionality includes easy customization of distance metrics, clustering algorithms, and SNF hyperparameters to generate diverse clustering solutions; calculation and plotting of associations between features, between patients, and between cluster solutions; and standard cluster validation approaches including resampled measures of cluster stability, standard metrics of cluster quality, and label propagation to evaluate generalizability in unseen data. Associated vignettes guide the user through using the package to identify patient subtypes while adhering to best practices for unsupervised learning.

**License** GPL (>= 3)

**Encoding** UTF-8

**RoxygenNote** 7.3.2

**Imports** cli, cluster, data.table, digest, dplyr, ggplot2, grDevices, MASS, mclust, methods, progressr, purrr, RColorBrewer, rlang, SNFtool, stats, tibble, tidyr, utils

**Suggests** circlize, ComplexHeatmap, InteractiveComplexHeatmap, clv, future, future.apply, knitr, rmarkdown, testthat (>= 3.0.0), ggalluvial, lifecycle, dbscan

**Config/testthat/edition** 3

**Depends** R (>= 4.1.0)

**LazyData** true

**VignetteBuilder** knitr

**URL** <https://branchlab.github.io/metasnf/>,  
<https://github.com/BRANCHlab/metasnf/>

**BugReports** <https://github.com/BRANCHlab/metasnf/issues>

**NeedsCompilation** no

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**Repository** CRAN

**Date/Publication** 2025-02-04 21:00:02 UTC

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abcd_anxiety	<i>Mock ABCD anxiety data</i>
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**Description**

A randomly shuffled and anonymized copy of anxiety data from the NIMH Data archive. The original file used was pdem02.txt. The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function `get_cbcl_anxiety`.

**Usage**

```
abcd_anxiety
```

**Format**

```
abcd_anxiety:
```

A data frame with 275 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset

**cbcl\_anxiety\_r** Ordinal value of impairment on CBCL anxiety, either 0 (no impairment), 1 (borderline clinical), or 2 (clinically impaired)

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

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abcd\_colour

*Mock ABCD "colour" data*

---

## Description

A randomly shuffled and anonymized copy of depression data from the NIMH Data archive. The original file used was pdem02.txt. The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function `get_cbc1_depress`. The data was transformed into categorical colour values to demonstrate the Chi-squared test capabilities of `extend_solutions`.

## Usage

```
abcd_colour
```

## Format

`abcd_colour`:

A data frame with 275 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset

**colour** Categorical transformation of `cbc1_depress`.

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and

follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

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abcd\_cort\_sa

*Mock ABCD cortical surface area data*


---

## Description

A randomly shuffled and anonymized copy of cortical surface area data from the NIMH Data archive. The original file used was `mrisdp10201.txt`. The file was pre-processed by the `abcdutils` package (<https://github.com/BRANCHlab/abcdutils>) function `get_cort_t`.

## Usage

```
abcd_cort_sa
```

## Format

```
abcd_cort_sa:
```

A data frame with 188 rows and 152 columns:

**patient** The unique identifier of the ABCD dataset

... Cortical surface areas of various ROIs (mm<sup>2</sup>, I think)

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete

listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

abcd\_cort\_t

*Mock ABCD cortical thickness data*

---

### Description

A randomly shuffled and anonymized copy of cortical thickness data from the NIMH Data archive. The original file used was `mrisd10201.txt`. The file was pre-processed by the `abcdutils` package (<https://github.com/BRANCHlab/abcdutils>) function `get_cort_t`.

### Usage

```
abcd_cort_t
```

### Format

```
abcd_cort_t:
```

A data frame with 188 rows and 152 columns:

**patient** The unique identifier of the ABCD dataset

... Cortical thicknesses of various ROIs (mm<sup>3</sup>, I think)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.



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`abcd_depress`*Mock ABCD depression data*

---

## Description

A randomly shuffled and anonymized copy of depression data from the NIMH Data archive. The original file used was pdem02.txt. The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function `get_cbc1_depress`.

## Usage

```
abcd_depress
```

## Format

`abcd_depress`:

A data frame with 275 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset

**cbc1\_depress\_r** Ordinal value of impairment on CBCL anxiety, either 0 (no impairment), 1 (borderline clinical), or 2 (clinically impaired)

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

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abcd_h_income	<i>Mock ABCD income data</i>
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### Description

Like abcd\_income, but with no NAs in patient column

### Usage

abcd\_h\_income

### Format

abcd\_income:

A data frame with 300 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset

**household\_income** Household income in 3 category levels (low = 1, medium = 2, high = 3)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

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abcd_income	<i>Mock ABCD income data</i>
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### Description

A randomly shuffled and anonymized copy of income data from the NIMH Data archive. The original file used was pdem02.txt The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function get\_income.

**Usage**

abcd\_income

**Format**

abcd\_income:

A data frame with 300 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset**household\_income** Household income in 3 category levels (low = 1, medium = 2, high = 3)**Source**

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

abcd\_pubertal

*Mock ABCD pubertal status data***Description**

A randomly shuffled and anonymized copy of pubertal status data from the NIMH Data archive. The original files used were abcd\_ssphp01.txt and abcd\_ssphy01.txt. The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function `get_pubertal_status`.

**Usage**

abcd\_pubertal

**Format**

abcd\_pubertal:

A data frame with 275 rows and 2 columns:

**patient** The unique identifier of the ABCD dataset**pubertal\_status** Average reported pubertal status between child and parent (1-5 categorical scale)

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

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abcd\_subc\_v

*Mock ABCD subcortical volumes data*

---

## Description

A randomly shuffled and anonymized copy of subcortical volume data from the NIMH Data archive. The original file used was smrip10201.txt The file was pre-processed by the abcdutils package (<https://github.com/BRANCHlab/abcdutils>) function `get_subc_v`.

## Usage

abcd\_subc\_v

## Format

abcd\_subc\_v:

A data frame with 174 rows and 31 columns:

**patient** The unique identifier of the ABCD dataset

... Subcortical volumes of various ROIs (mm<sup>3</sup>, I think)

## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and

follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

add\_settings\_df\_rows *Add rows to a settings\_df*

---

## Description

Add rows to a settings\_df

## Usage

```
add_settings_df_rows(
  sdf,
  n_solutions = 0,
  min_removed_inputs = 0,
  max_removed_inputs = sum(startsWith(colnames(sdf), "inc_")) - 1,
  dropout_dist = "exponential",
  min_alpha = NULL,
  max_alpha = NULL,
  min_k = NULL,
  max_k = NULL,
  min_t = NULL,
  max_t = NULL,
  alpha_values = NULL,
  k_values = NULL,
  t_values = NULL,
  possible_snf_schemes = c(1, 2, 3),
  clustering_algorithms = NULL,
  continuous_distances = NULL,
  discrete_distances = NULL,
  ordinal_distances = NULL,
  categorical_distances = NULL,
  mixed_distances = NULL,
  dfl = NULL,
  snf_input_weights = NULL,
  snf_domain_weights = NULL,
  retry_limit = 10,
```

```

    allow_duplicates = FALSE
  )

```

### Arguments

sdf	The existing settings data frame
n_solutions	Number of rows to generate for the settings data frame.
min_removed_inputs	The smallest number of input data frames that may be randomly removed. By default, 0.
max_removed_inputs	The largest number of input data frames that may be randomly removed. By default, this is 1 less than all the provided input data frames in the data list.
dropout_dist	Parameter controlling how the random removal of input data frames should occur. Can be "none" (no input data frames are randomly removed), "uniform" (uniformly sample between min_removed_inputs and max_removed_inputs to determine number of input data frames to remove), or "exponential" (pick number of input data frames to remove by sampling from min_removed_inputs to max_removed_inputs with an exponential distribution; the default).
min_alpha	The minimum value that the alpha hyperparameter can have. Random assigned value of alpha for each row will be obtained by uniformly sampling numbers between min_alpha and max_alpha at intervals of 0.1. Cannot be used in conjunction with the alpha_values parameter.
max_alpha	The maximum value that the alpha hyperparameter can have. See min_alpha parameter. Cannot be used in conjunction with the alpha_values parameter.
min_k	The minimum value that the k hyperparameter can have. Random assigned value of k for each row will be obtained by uniformly sampling numbers between min_k and max_k at intervals of 1. Cannot be used in conjunction with the k_values parameter.
max_k	The maximum value that the k hyperparameter can have. See min_k parameter. Cannot be used in conjunction with the k_values parameter.
min_t	The minimum value that the t hyperparameter can have. Random assigned value of t for each row will be obtained by uniformly sampling numbers between min_t and max_t at intervals of 1. Cannot be used in conjunction with the t_values parameter.
max_t	The maximum value that the t hyperparameter can have. See min_t parameter. Cannot be used in conjunction with the t_values parameter.
alpha_values	A number or numeric vector of a set of possible values that alpha can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_alpha or max_alpha parameters.
k_values	A number or numeric vector of a set of possible values that k can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_k or max_k parameters.
t_values	A number or numeric vector of a set of possible values that t can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_t or max_t parameters.

possible_snf_schemes	A vector containing the possible snf_schemes to uniformly randomly select from. By default, the vector contains all 3 possible schemes: c(1, 2, 3). 1 corresponds to the "individual" scheme, 2 corresponds to the "domain" scheme, and 3 corresponds to the "twostep" scheme.
clustering_algorithms	A list of clustering algorithms to uniformly randomly pick from when clustering. When not specified, randomly select between spectral clustering using the eigen-gap heuristic and spectral clustering using the rotation cost heuristic. See ?clust_fns_list for more details on running custom clustering algorithms.
continuous_distances	A vector of continuous distance metrics to use when a custom dist_fns_list is provided.
discrete_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
ordinal_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
categorical_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
mixed_distances	A vector of mixed distance metrics to use when a custom dist_fns_list is provided.
df1	List containing distance metrics to vary over. See ?generate_dist_fns_list.
snf_input_weights	Nested list containing weights for when SNF is used to merge individual input measures (see ?generate_snf_weights)
snf_domain_weights	Nested list containing weights for when SNF is used to merge domains (see ?generate_snf_weights)
retry_limit	The maximum number of attempts to generate a novel row. This function does not return matrices with identical rows. As the range of requested possible settings tightens and the number of requested rows increases, the risk of randomly generating a row that already exists increases. If a new random row has matched an existing row retry_limit number of times, the function will terminate.
allow_duplicates	If TRUE, enables creation of a settings data frame with duplicate non-feature weighting related hyperparameters. This function should only be used when paired with a custom weights matrix that has non-duplicate rows.

**Value**

A settings data frame

---

adjusted\_rand\_index\_heatmap

*Heatmap of pairwise adjusted rand indices between solutions*


---

### Description

**[Deprecated]** Defunct function to create an ARI heatmap. Please use `meta_cluster_heatmap()` instead.

### Usage

```
adjusted_rand_index_heatmap(
  aris,
  order = NULL,
  cluster_rows = FALSE,
  cluster_columns = FALSE,
  log_graph = FALSE,
  scale_diag = "none",
  min_colour = "#282828",
  max_colour = "firebrick2",
  col = circlize::colorRamp2(c(min(aris), max(aris)), c(min_colour, max_colour)),
  ...
)
```

### Arguments

<code>aris</code>	Matrix of adjusted rand indices from <code>calc_aris()</code>
<code>order</code>	Numeric vector containing row order of the heatmap.
<code>cluster_rows</code>	Whether rows should be clustered.
<code>cluster_columns</code>	Whether columns should be clustered.
<code>log_graph</code>	If TRUE, log transforms the graph.
<code>scale_diag</code>	Method of rescaling matrix diagonals. Can be "none" (don't change diagonals), "mean" (replace diagonals with average value of off-diagonals), or "zero" (replace diagonals with 0).
<code>min_colour</code>	Colour used for the lowest value in the heatmap.
<code>max_colour</code>	Colour used for the highest value in the heatmap.
<code>col</code>	Colour ramp to use for the heatmap.
<code>...</code>	Additional parameters passed to <code>similarity_matrix_heatmap()</code> , the function that this function wraps.

### Value

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the pairwise adjusted Rand indices (similarities) between the cluster solutions of the provided solutions data frame.



---

age_df	<i>Mock age data</i>
--------	----------------------

---

**Description**

Mock age data

**Usage**

```
age_df
```

**Format**

age\_df:

A data frame with 200 rows and 2 columns:

**patient\_id** Random three-digit number uniquely identifying the patient

**age** Mock age feature

**Source**

This data came from the SNFtool package, with slight modifications.

---

alluvial_cluster_plot	<i>Alluvial plot of patients across cluster counts and important features</i>
-----------------------	---

---

**Description**

This alluvial plot shows how observations in a similarity matrix could have been clustered over a set of clustering functions.

**Usage**

```
alluvial_cluster_plot(  
  cluster_sequence,  
  similarity_matrix,  
  dl = NULL,  
  data = NULL,  
  key_outcome,  
  key_label = key_outcome,  
  extra_outcomes = NULL,  
  title = NULL  
)
```

**Arguments**

<code>cluster_sequence</code>	A list of clustering algorithms.
<code>similarity_matrix</code>	A similarity matrix.
<code>dl</code>	A data list.
<code>data</code>	A data frame that contains any features to include in the plot.
<code>key_outcome</code>	The name of the feature that determines how each patient stream is coloured in the alluvial plot.
<code>key_label</code>	Name of key outcome to be used for the plot legend.
<code>extra_outcomes</code>	Names of additional features to add to the plot.
<code>title</code>	Title of the plot.

**Value**

An alluvial plot (class "gg" and "ggplot") showing distribution of a feature across varying number cluster solutions.

**Examples**

```
input_dl <- data_list(
  list(gender_df, "gender", "demographics", "categorical"),
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),
  uid = "patient_id"
)

sc <- snf_config(input_dl, n_solutions = 1)

sol_df <- batch_snf(input_dl, sc, return_sim_mats = TRUE)

sim_mats <- sim_mats_list(sol_df)

clust_fn_sequence <- list(spectral_two, spectral_four)

alluvial_cluster_plot(
  cluster_sequence = clust_fn_sequence,
  similarity_matrix = sim_mats[[1]],
  dl = input_dl,
  key_outcome = "gender", # the name of the feature of interest
  key_label = "Gender", # how the feature of interest should be displayed
  extra_outcomes = "diagnosis", # more features to plot but not colour by
  title = "Gender Across Cluster Counts"
)
```

---

anxiety	<i>Mock ABCD anxiety data</i>
---------	-------------------------------

---

### Description

Like the mock data frame "abcd\_colour", but with "unique\_id" as the "uid".

### Usage

```
anxiety
```

### Format

```
anxiety:
```

A data frame with 275 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**cbcl\_anxiety\_r** Ordinal value of impairment on CBCL anxiety, either 0 (no impairment), 1 (borderline clinical), or 2 (clinically impaired)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

arrange	<i>Arrange rows in an object</i>
---------	----------------------------------

---

**Description**

Arrange rows in an object

**Usage**

```
arrange(.data, ...)
```

**Arguments**

.data	The object to arrange columns from.
...	Additional arguments for arranging.

**Value**

Object with arranged columns.

---

as.data.frame.data_list	<i>Coerce a data_list class object into a data.frame class object</i>
-------------------------	---

---

**Description**

Horizontally joins data frames within a data list into a single data frame, using the uid attribute as the joining key.

**Usage**

```
## S3 method for class 'data_list'
as.data.frame(x, row.names = NULL, optional = FALSE, ...)
```

**Arguments**

x	A data_list class object.
row.names	Additional parameter passed to as.data.frame().
optional	Additional parameter passed to as.data.frame().
...	Additional parameter passed to as.data.frame().

**Value**

dl\_df A data.frame class object with all the features and observations of dl.

---

```
as.data.frame.ext_solutions_df
    Coerce a ext_solutions_df class object into a data.frame class
    object
```

---

**Description**

Coerce a ext\_solutions\_df class object into a data.frame class object

**Usage**

```
## S3 method for class 'ext_solutions_df'
as.data.frame(
  x,
  row.names = NULL,
  optional = FALSE,
  keep_attributes = FALSE,
  ...
)
```

**Arguments**

x	A ext_solutions_df class object.
row.names	Additional parameter passed to as.data.frame().
optional	Additional parameter passed to as.data.frame().
keep_attributes	If TRUE, resulting data frame includes settings data frame and weights matrix.
...	Additional parameter passed to as.data.frame().

**Value**

A data.frame class object with all the columns of x and its contained solutions data frame.

---

```
as.data.frame.solutions_df
    Coerce a solutions_df class object into a data.frame class object
```

---

**Description**

Coerce a solutions\_df class object into a data.frame class object

**Usage**

```
## S3 method for class 'solutions_df'
as.data.frame(
  x,
  row.names = NULL,
  optional = FALSE,
  keep_attributes = FALSE,
  ...
)
```

**Arguments**

x	A solutions_df class object.
row.names	Additional parameter passed to as.data.frame().
optional	Additional parameter passed to as.data.frame().
keep_attributes	If TRUE, resulting data frame includes settings data frame and weights matrix.
...	Additional parameter passed to as.data.frame().

**Value**

A data.frame class object with all the columns of x and its contained solutions data frame.

---

assemble_data	<i>Collapse a data frame and/or a data list into a single data frame</i>
---------------	--

---

**Description**

Collapse a data frame and/or a data list into a single data frame

**Usage**

```
assemble_data(data, dl)
```

**Arguments**

data	A data frame.
dl	A nested list of input data from data_list().

**Value**

A class "data.frame" object containing all the features of the provided data frame and/or data list.

---

assoc\_pval\_heatmap      *Heatmap of pairwise associations between features*

---

### Description

Heatmap of pairwise associations between features

### Usage

```
assoc_pval_heatmap(
  correlation_matrix,
  scale_diag = "max",
  cluster_rows = TRUE,
  cluster_columns = TRUE,
  show_row_names = TRUE,
  show_column_names = TRUE,
  show_heatmap_legend = FALSE,
  confounders = NULL,
  out_of_models = NULL,
  annotation_colours = NULL,
  labels_colour = NULL,
  split_by_domain = FALSE,
  dl = NULL,
  significance_stars = TRUE,
  slice_font_size = 8,
  ...
)
```

### Arguments

correlation_matrix	Matrix containing all pairwise association p-values. The recommended way to obtain this matrix is through the <code>calc_assoc_pval</code> function.
scale_diag	Parameter that controls how the diagonals of the <code>correlation_matrix</code> are adjusted in the heatmap. For best viewing, this is set to "max", which will match the diagonals to whichever pairwise association has the highest p-value.
cluster_rows	Parameter for <code>ComplexHeatmap::Heatmap</code> . Will be ignored if <code>split_by_domain</code> is also provided.
cluster_columns	Parameter for <code>ComplexHeatmap::Heatmap</code> . Will be ignored if <code>split_by_domain</code> is also provided.
show_row_names	Parameter for <code>ComplexHeatmap::Heatmap</code> .
show_column_names	Parameter for <code>ComplexHeatmap::Heatmap</code> .
show_heatmap_legend	Parameter for <code>ComplexHeatmap::Heatmap</code> .

confounders	A named list where the elements are columns in the correlation_matrix and the names are the corresponding display names.
out_of_models	Like confounders, but a named list of out of model measures (who are also present as columns in the correlation_matrix).
annotation_colours	Named list of heatmap annotations and their colours.
labels_colour	Vector of colours to use for the columns and rows of the heatmap.
split_by_domain	The results of dl_var_summar - a data frame that has the domain of every feature in the plotted data. columns of the correlation_matrix. Will be used to "slice" the heatmap into visually separated sections.
dl	A nested list of input data from data_list().
significance_stars	If TRUE (default), plots significance stars on heatmap cells
slice_font_size	Font size for domain separating labels.
...	Additional parameters passed into ComplexHeatmap::Heatmap.

**Value**

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the pairwise associations between features from the provided correlation\_matrix.

**Examples**

```
#data_list <- data_list(
#  list(income, "household_income", "demographics", "ordinal"),
#  list(pubertal, "pubertal_status", "demographics", "continuous"),
#  list(fav_colour, "favourite_colour", "demographics", "categorical"),
#  list(anxiety, "anxiety", "behaviour", "ordinal"),
#  list(depress, "depressed", "behaviour", "ordinal"),
#  uid = "unique_id"
#)
#
#assoc_pval_matrix <- calc_assoc_pval_matrix(data_list)
#ap_heatmap <- assoc_pval_heatmap(assoc_pval_matrix)
```

---

as\_ari\_matrix

*Convert an object to an ARI matrix*


---

**Description**

This function coerces non-ari\_matrix class objects into ari\_matrix class objects.

**Usage**

```
as_ari_matrix(x)
```



**Arguments**

x                    The object to convert into a weights matrix.

**Value**

An `ari_matrix` class object.

---

as\_data\_list                    *Convert an object to a data list*

---

**Description**

This function coerces non-`data_list` class objects into `data_list` class objects.

**Usage**

```
as_data_list(x)
```

**Arguments**

x                    The object to convert into a data list.

**Value**

A `data_list` class object.

---

as\_settings\_df                    *Convert an object to a settings data frame*

---

**Description**

This function coerces non-`settings_df` class objects into `settings_df` class objects.

**Usage**

```
as_settings_df(x)
```

**Arguments**

x                    The object to convert into a data list.

**Value**

A `settings_df` class object.

---

as_sim_mats_list	<i>Convert an object to a similarity matrix list</i>
------------------	--

---

**Description**

This function converts non-sim\_mats\_list class objects into sim\_mats\_list class objects.

**Usage**

```
as_sim_mats_list(x)
```

**Arguments**

x                    The object to convert into a sim\_mats\_list. Must be a list of square matrices with identical column and row names.

**Value**

A sim\_mats\_list class object.

---

as_snf_config	<i>Convert an object to a snf config</i>
---------------	--

---

**Description**

This function coerces non-snf\_config class objects into snf\_config class objects.

**Usage**

```
as_snf_config(x)
```

**Arguments**

x                    The object to convert into a snf config.

**Value**

A snf\_config class object.

---

as_weights_matrix	<i>Convert an object to a weights matrix</i>
-------------------	--

---

**Description**

This function converts non-weights\_matrix objects into weights\_matrix class objects.

**Usage**

```
as_weights_matrix(x)
```

**Arguments**

x                    The object to convert into a data list.

**Value**

A weights\_matrix class object.

---

auto_plot	<i>Automatically plot features across clusters</i>
-----------	--

---

**Description**

Given a single row of a solutions data frame and data provided through a data list, this function will return a series of bar and/or jitter plots based on feature types.

**Usage**

```
auto_plot(  
  sol_df_row = NULL,  
  dl = NULL,  
  cluster_df = NULL,  
  return_plots = TRUE,  
  save = NULL,  
  jitter_width = 6,  
  jitter_height = 6,  
  bar_width = 6,  
  bar_height = 6,  
  verbose = FALSE  
)
```

**Arguments**

sol_df_row	A single row of a solutions data frame.
dl	A data list containing data to plot.
cluster_df	Directly provide a cluster_df rather than a solutions matrix. Useful if plotting data from label propagated results.
return_plots	If TRUE, the function will return a list of plots. If FALSE, the function will instead return the full data frame used for plotting.
save	If a string is provided, plots will be saved and this string will be used to prefix plot names.
jitter_width	Width of jitter plots if save is specified.
jitter_height	Height of jitter plots if save is specified.
bar_width	Width of bar plots if save is specified.
bar_height	Height of bar plots if save is specified.
verbose	If TRUE, output progress to console.

**Value**

By default, returns a list of plots (class "gg", "ggplot") with one plot for every feature in the provided data list and/or target list. If return\_plots is FALSE, will instead return a single "data.frame" object containing every provided feature for every observation in long format.

---

bar_plot	<i>Bar plot separating a feature by cluster</i>
----------	---

---

**Description**

Bar plot separating a feature by cluster

**Usage**

```
bar_plot(df, feature)
```

**Arguments**

df	A data.frame containing cluster column and the feature to plot.
feature	The feature to plot.

**Value**

A bar plot (class "gg", "ggplot") showing the distribution of a feature across clusters.

---

batch_snf	<i>Run variations of SNF.</i>
-----------	-------------------------------

---

## Description

This is the core function of the `metasnf` package. Using the information stored in a `settings_df` (see `?settings_df`) and a data list (see `?data_list`), run repeated complete SNF pipelines to generate a broad space of post-SNF cluster solutions.

## Usage

```
batch_snf(dl, sc, processes = 1, return_sim_mats = FALSE, sim_mats_dir = NULL)
```

## Arguments

<code>dl</code>	A nested list of input data from <code>data_list()</code> .
<code>sc</code>	An <code>snf_config</code> class object which stores all sets of hyperparameters used to transform data in <code>dl</code> into a cluster solutions. See <code>?settings_df</code> or <a href="https://branchlab.github.io/metasnf/arti">https://branchlab.github.io/metasnf/arti</a> for more details.
<code>processes</code>	Specify number of processes used to complete SNF iterations <ul style="list-style-type: none"> <li>• 1 (default) Sequential processing: function will iterate through the <code>settings_df</code> one row at a time with a for loop. This option will not make use of multiple CPU cores, but will show a progress bar.</li> <li>• 2 or higher: Parallel processing will use the <code>future.apply::future_apply</code> to distribute the SNF iterations across the specified number of CPU cores. If higher than the number of available cores, a warning will be raised and the maximum number of cores will be used.</li> <li>• max: All available cores will be used.</li> </ul>
<code>return_sim_mats</code>	If TRUE, function will return a list where the first element is the solutions data frame and the second element is a list of similarity matrices for each row in the <code>sol_df</code> . Default FALSE.
<code>sim_mats_dir</code>	If specified, this directory will be used to save all generated similarity matrices.

## Value

By default, returns a solutions data frame (class "data.frame"), a a data frame containing one row for every row of the provided settings matrix, all the original columns of that settings data frame, and new columns containing the assigned cluster of each observation from the cluster solution derived by that row's settings. If `return_sim_mats` is TRUE, the function will instead return a list containing the solutions data frame as well as a list of the final similarity matrices (class "matrix") generated by SNF for each row of the settings data frame. If `suppress_clustering` is TRUE, the solutions data frame will not be returned in the output.

**Examples**

```

input_dl <- data_list(
  list(gender_df, "gender", "demographics", "categorical"),
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),
  uid = "patient_id"
)

sc <- snf_config(input_dl, n_solutions = 3)

# A solutions data frame without similarity matrices:
sol_df <- batch_snf(input_dl, sc)

# A solutions data frame with similarity matrices:
# sol_df <- batch_snf(input_dl, sc, return_sim_mats = TRUE)
# sim_mats_list(sol_df)

```

---

batch\_snf\_subsamples *Run SNF clustering pipeline on a list of subsampled data lists.*

---

**Description**

Run SNF clustering pipeline on a list of subsampled data lists.

**Usage**

```

batch_snf_subsamples(
  dl_subsamples,
  sc,
  processes = 1,
  return_sim_mats = FALSE,
  sim_mats_dir = NULL,
  verbose = TRUE
)

```

**Arguments**

- |               |  |
|---------------|--|
| dl_subsamples | A list of subsampled data lists. This object is generated by the function <code>batch_snf_subsamples()</code> .  |
| sc            | An <code>snf_config</code> class object which stores all sets of hyperparameters used to transform data in <code>dl</code> into a cluster solutions. See <code>?settings_df</code> or <a href="https://branchlab.github.io/metasnfn/articles/">https://branchlab.github.io/metasnfn/articles/</a> for more details.  |
| processes     | Specify number of processes used to complete SNF iterations <ul style="list-style-type: none"> <li>• 1 (default) Sequential processing: function will iterate through the <code>settings_df</code> one row at a time with a for loop. This option will not make use of multiple CPU cores, but will show a progress bar.</li> <li>• 2 or higher: Parallel processing will use the <code>future.apply::future_apply</code> to distribute the SNF iterations across the specified number of CPU cores. If higher than the number of available cores, a warning will be raised and the maximum number of cores will be used.</li> </ul> |

- max: All available cores will be used.
- return\_sim\_mats      If TRUE, function will return a list where the first element is the solutions data frame and the second element is a list of similarity matrices for each row in the sol\_df. Default FALSE.
- sim\_mats\_dir        If specified, this directory will be used to save all generated similarity matrices.
- verbose             If TRUE, output progress to console.

### Value

By default, returns a one-element list: cluster\_solutions, which is itself a list of cluster solution data frames corresponding to each of the provided data list subsamples. Setting the parameters return\_sim\_mats and return\_solutions to TRUE will turn the result of the function to a three-element list containing the corresponding solutions data frames and final fused similarity matrices of those cluster solutions, should you require these objects for your own stability calculations.

### Examples

```
# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(my_dl, n_solutions = 5, max_k = 40)
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
#
# batch_subsample_results <- batch_snf_subsamples(
#   my_dl_subsamples,
#   sc,
#   verbose = TRUE
# )
```

---

calculate\_coclustering

*Calculate coclustering data.*

---

### Description

Calculate coclustering data.

**Usage**

```
calculate_coclustering(subsample_solutions, sol_df, verbose = FALSE)
```

**Arguments**

`subsample_solutions` A list of containing cluster solutions from distinct subsamples of the data. This object is generated by the function `batch_snf_subsamples()`. These solutions should correspond to the ones in the solutions data frame.

`sol_df` A solutions data frame. This object is generated by the function `batch_snf()`. The solutions in the solutions data frame should correspond to those in the subsample solutions.

`verbose` If TRUE, output time remaining estimates to console.

**Value**

A list containing the following components:

- `cocluster_dfs`: A list of data frames, one per cluster solution, that shows the number of times that every pair of observations in the original cluster solution occurred in the same subsample, the number of times that every pair clustered together in a subsample, and the corresponding fraction of times that every pair clustered together in a subsample.
- `cocluster_ss_mats`: The number of times every pair of observations occurred in the same subsample, formatted as a pairwise matrix.
- `cocluster_sc_mats`: The number of times every pair of observations occurred in the same cluster, formatted as a pairwise matrix.
- `cocluster_cf_mats`: The fraction of times every pair of observations occurred in the same cluster, formatted as a pairwise matrix.
- `cocluster_summary`: Specifically among pairs of observations that clustered together in the original full cluster solution, what fraction of those pairs remained clustered together throughout the subsample solutions. This information is formatted as a data frame with one row per cluster solution.

**Examples**

```
# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(my_dl, n_solutions = 5, max_k = 40)
#
# sol_df <- batch_snf(my_dl, sc)
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
```



```

#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
#
# batch_subsample_results <- batch_snf_subsamples(
#   my_dl_subsamples,
#   sc,
#   verbose = TRUE
# )
#
# coclustering_results <- calculate_coclustering(
#   batch_subsample_results,
#   sol_df,
#   verbose = TRUE
# )

```

---

calc\_aris

*Construct an ARI matrix storing inter-solution similarities*


---

### Description

This function constructs an `ari_matrix` class object from a `solutions_df` class object. The ARI matrix stores pairwise adjusted Rand indices for all cluster solutions as well as a numeric order for the solutions data frame based on the hierarchical clustering of the ARI matrix.

### Usage

```

calc_aris(
  sol_df,
  processes = 1,
  verbose = FALSE,
  dist_method = "euclidean",
  hclust_method = "complete"
)

```

### Arguments

<code>sol_df</code>	Solutions data frame containing cluster solutions to calculate pairwise ARIs for.
<code>processes</code>	Specify number of processes used to complete calculations <ul style="list-style-type: none"> <li>• 1 (default) Sequential processing</li> <li>• 2 or higher: Parallel processing will use the <code>future::future_apply</code> to distribute the calculations across the specified number of CPU cores. If higher than the number of available cores, a warning will be raised and the maximum number of cores will be used.</li> <li>• max: All available cores will be used. Note that no progress indicator is available during multi-core processing.</li> </ul>
<code>verbose</code>	If TRUE, output progress to console.

`dist_method` Distance method to use when calculating sorting order to of the matrix. Argument is directly passed into `stats::dist`. Options include "euclidean", "maximum", "manhattan", "canberra", "binary", or "minkowski".

`hclust_method` Agglomerative method to use when calculating sorting order by `stats::hclust`. Options include "ward.D", "ward.D2", "single", "complete", "average", "mcquitty", "median", or "centroid".

**Value**

om\_aris ARIs between clustering solutions of an solutions data frame

**Examples**

```
d1 <- data_list(
  list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  uid = "unique_id"
)

sc <- snf_config(d1, n_solutions = 3)
sol_df <- batch_snf(d1, sc)
calc_aris(sol_df)
```

---

calc\_assoc\_pval\_matrix

*Calculate p-values for all pairwise associations of features in a data list*

---

**Description**

Calculate p-values for all pairwise associations of features in a data list

**Usage**

```
calc_assoc_pval_matrix(d1, verbose = FALSE, cat_test = "chi_squared")
```

**Arguments**

`d1` A nested list of input data from `data_list()`.

`verbose` If TRUE, output progress to the console.

`cat_test` String indicating which statistical test will be used to associate cluster with a categorical feature. Options are "chi\_squared" for the Chi-squared test and "fisher\_exact" for Fisher's exact test.

**Value**

A "matrix" class object containing pairwise association p-values between the features in the provided data list.

**Examples**

```

data_list <- data_list(
  list(income, "household_income", "demographics", "ordinal"),
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  list(anxiety, "anxiety", "behaviour", "ordinal"),
  list(depress, "depressed", "behaviour", "ordinal"),
  uid = "unique_id"
)

assoc_pval_matrix <- calc_assoc_pval_matrix(data_list)

```

calc\_nmis

*Calculate feature NMIs for a data list and a solutions data frame***Description**

Normalized mutual information scores can be used to indirectly measure how important a feature may have been in producing a cluster solution. This function will calculate the normalized mutual information between cluster solutions in a solutions data frame as well as cluster solutions created by including only a single feature from a provided data list, but otherwise using all the same hyperparameters as specified in the original SNF config. Note that NMIs can be calculated between two cluster solutions regardless of what features were actually used to create those cluster solutions. For example, a feature that was not involved in producing a particular cluster solution may still have a high NMI with that cluster solution (typically because it was highly correlated with a different feature that was used).

**Usage**

```

calc_nmis(
  dl,
  sol_df,
  transpose = TRUE,
  ignore_inclusions = TRUE,
  verbose = FALSE
)

```

**Arguments**

dl	A nested list of input data from <code>data_list()</code> .
sol_df	Result of <code>batch_snf</code> storing cluster solutions and the settings that were used to generate them. Use the same value as was used in the original call to <code>batch_snf()</code> .
transpose	If TRUE, will transpose the output data frame.
ignore_inclusions	If TRUE, will ignore the inclusion columns in the solutions data frame and calculate NMIs for all features. If FALSE, will give NAs for features that were dropped on a given <code>settings_df</code> row.
verbose	If TRUE, output progress to console.

**Value**

A "data.frame" class object containing one row for every feature in the provided data list and one column for every solution in the provided solutions data frame. Populated values show the calculated NMI score for each feature-solution combination.

**Examples**

```
input_dl <- data_list(  
  list(gender_df, "gender", "demographics", "categorical"),  
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),  
  uid = "patient_id"  
)  
  
sc <- snf_config(input_dl, n_solutions = 2)  
  
sol_df <- batch_snf(input_dl, sc)  
  
calc_nmis(input_dl, sol_df)
```

---

cancer\_diagnosis\_df    *Mock diagnosis data*

---

**Description**

This is the same data as `diagnosis_df`, with renamed features and columns.

**Usage**

```
cancer_diagnosis_df
```

**Format**

`cancer_diagnosis_df`:

A data frame with 200 rows and 2 columns:

**patient\_id** Random three-digit number uniquely identifying the patient

**diagnosis** Mock cancer diagnosis feature (1, 2, or 3)

**Source**

This data came from the SNFtool package, with slight modifications.

---

cell\_significance\_fn *Place significance stars on ComplexHeatmap cells.*

---

**Description**

This is an internal function meant to be used to by the assoc\_pval\_heatmap function.

**Usage**

```
cell_significance_fn(data)
```

**Arguments**

data                    The matrix containing the cells to base the significance stars on.

**Value**

cell\_fn Another function that is well-formatted for usage as the cell\_fun argument in ComplexHeatmap::Heatmap.

---

check\_dataless\_annotations

*Helper function to stop annotation building when no data was provided*

---

**Description**

Helper function to stop annotation building when no data was provided

**Usage**

```
check_dataless_annotations(annotation_requests, data)
```

**Arguments**

annotation\_requests

A list of requested annotations

data

A data frame with data to build annotations

**Value**

Does not return any value. This function just raises an error when annotations are requested without any provided data for a heatmap.

check\_hm\_dependencies *Check for ComplexHeatmap and circlize dependencies*

---

**Description**

Check for ComplexHeatmap and circlize dependencies

**Usage**

```
check_hm_dependencies()
```

**Value**

Does not return any value. This function just checks that the ComplexHeatmap and circlize packages are installed.

---

check\_similarity\_matrices  
*Check validity of similarity matrices*

---

**Description**

Check to see if similarity matrices in a list have the following properties:

1. The maximum value in the entire matrix is 0.5
2. Every value in the diagonal is 0.5

**Usage**

```
check_similarity_matrices(similarity_matrices)
```

**Arguments**

similarity\_matrices  
A list of similarity matrices

**Value**

valid\_matrices Boolean indicating if properties are met by all similarity matrices

---

`clust_fns`*Built-in clustering algorithms*

---

**Description**

These functions can be used when building a `metasnf` clustering functions list. Each function converts a similarity matrix (matrix class object) to a cluster solution (numeric vector). Note that these functions (or custom clustering functions) cannot accept number of clusters as a parameter; this value must be built into the function itself if necessary.

**Usage**

```
spectral_eigen(similarity_matrix)
spectral_rot(similarity_matrix)
spectral_eigen_classic(similarity_matrix)
spectral_rot_classic(similarity_matrix)
spectral_two(similarity_matrix)
spectral_three(similarity_matrix)
spectral_four(similarity_matrix)
spectral_five(similarity_matrix)
spectral_six(similarity_matrix)
spectral_seven(similarity_matrix)
spectral_eight(similarity_matrix)
spectral_nine(similarity_matrix)
spectral_ten(similarity_matrix)
```

**Arguments**

```
similarity_matrix
    A similarity matrix.
```

**Details**

- `spectral_eigen`: Spectral clustering where the number of clusters is based on the eigen-gap heuristic

- spectral\_rot: Spectral clustering where the number of clusters is based on the rotation-cost heuristic
- spectral\_(C): Spectral clustering for a C-cluster solution.

### Value

solution\_data A vector of cluster assignments

---

clust\_fns\_list            *Build a clustering algorithms list*

---

### Description

This function can be used to specify custom clustering algorithms to apply to the final similarity matrices produced by each run of the batch\_snf function.

### Usage

```
clust_fns_list(clust_fns = NULL, use_default_clust_fns = FALSE)
```

### Arguments

clust\_fns            A list of named clustering functions

use\_default\_clust\_fns  
If TRUE, prepend the base clustering algorithms (spectral\_eigen and spectral\_rot, which apply spectral clustering and use the eigen-gap and rotation cost heuristics respectively for determining the number of clusters in the graph) to clust\_fns.

### Value

A list of clustering algorithm functions that can be passed into the batch\_snf and generate\_settings\_list functions.

### Examples

```
# Using just the base clustering algorithms -----
# This will just contain spectral_eigen and spectral_rot
cfl <- clust_fns_list(use_default_clust_fns = TRUE)

# Adding algorithms provided by the package -----
# This will contain the base clustering algorithms (spectral_eigen,
# spectral_rot) as well as two pre-defined spectral clustering functions
# that force the number of clusters to be two or five
cfl <- clust_fns_list(
  clust_fns = list(
    "two_cluster_spectral" = spectral_two,
    "five_cluster_spectral" = spectral_five
  )
)
```



```

)

# Adding your own algorithms -----
# This will contain the base and user-provided clustering algorithms
my_clustering_algorithm <- function(similarity_matrix) {
  # your code that converts similarity matrix to clusters here...
}

# Suppress the base algorithms-----
# This will contain only user-provided clustering algorithms
cfl <- clust_fns_list(
  clust_fns = list(
    "two_cluster_spectral" = spectral_two,
    "five_cluster_spectral" = spectral_five
  )
)

```

---

cocluster\_density      *Density plot coclustering stability across subsampled data.*

---

### Description

This function creates a density plot that shows, for all pairs of observations that originally clustered together, the distribution of the the fractions that those pairs clustered together across subsampled data.

### Usage

```
cocluster_density(cocluster_df)
```

### Arguments

`cocluster_df`      A data frame containing coclustering data for a single cluster solution. This object is generated by the `calculate_coclustering` function.

### Value

Density plot (class "gg", "ggplot") of the distribution of coclustering across pairs and subsamples of the data.

### Examples

```

# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
# )
#

```

```

# sc <- snf_config(my_dl, n_solutions = 5, max_k = 40)
#
# sol_df <- batch_snf(my_dl, sc)
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
#
# batch_subsample_results <- batch_snf_subsamples(
#   my_dl_subsamples,
#   sc,
#   verbose = TRUE
# )
#
# coclustering_results <- calculate_coclustering(
#   batch_subsample_results,
#   sol_df,
#   verbose = TRUE
# )
#
# cocluster_density(cocluster_dfs[[1]])

```

---

cocluster\_heatmap

*Heatmap of observation co-clustering across resampled data.*


---

### Description

Create a heatmap that shows the distribution of observation co-clustering across resampled data.

### Usage

```

cocluster_heatmap(
  cocluster_df,
  cluster_rows = TRUE,
  cluster_columns = TRUE,
  show_row_names = FALSE,
  show_column_names = FALSE,
  dl = NULL,
  data = NULL,
  left_bar = NULL,
  right_bar = NULL,
  top_bar = NULL,
  bottom_bar = NULL,
  left_hm = NULL,
  right_hm = NULL,
  top_hm = NULL,
  bottom_hm = NULL,

```

```

    annotation_colours = NULL,
    min_colour = NULL,
    max_colour = NULL,
    ...
)

```

### Arguments

<code>cocluster_df</code>	A data frame containing coclustering data for a single cluster solution. This object is generated by the <code>calculate_coclustering</code> function.
<code>cluster_rows</code>	Argument passed to <code>ComplexHeatmap::Heatmap()</code> .
<code>cluster_columns</code>	Argument passed to <code>ComplexHeatmap::Heatmap()</code> .
<code>show_row_names</code>	Argument passed to <code>ComplexHeatmap::Heatmap()</code> .
<code>show_column_names</code>	Argument passed to <code>ComplexHeatmap::Heatmap()</code> .
<code>dl</code>	See <code>?similarity_matrix_heatmap</code> .
<code>data</code>	See <code>?similarity_matrix_heatmap</code> .
<code>left_bar</code>	See <code>?similarity_matrix_heatmap</code> .
<code>right_bar</code>	See <code>?similarity_matrix_heatmap</code> .
<code>top_bar</code>	See <code>?similarity_matrix_heatmap</code> .
<code>bottom_bar</code>	See <code>?similarity_matrix_heatmap</code> .
<code>left_hm</code>	See <code>?similarity_matrix_heatmap</code> .
<code>right_hm</code>	See <code>?similarity_matrix_heatmap</code> .
<code>top_hm</code>	See <code>?similarity_matrix_heatmap</code> .
<code>bottom_hm</code>	See <code>?similarity_matrix_heatmap</code> .
<code>annotation_colours</code>	See <code>?similarity_matrix_heatmap</code> .
<code>min_colour</code>	See <code>?similarity_matrix_heatmap</code> .
<code>max_colour</code>	See <code>?similarity_matrix_heatmap</code> .
<code>...</code>	Arguments passed to <code>ComplexHeatmap::Heatmap()</code> .

### Value

Heatmap (class "Heatmap" from `ComplexHeatmap`) object showing the distribution of observation co-clustering across resampled data.

### Examples

```

# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"

```

```

# )
#
# sc <- snf_config(my_dl, n_solutions = 5, max_k = 40)
#
# sol_df <- batch_snf(my_dl, sc)
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
#
# batch_subsample_results <- batch_snf_subsamples(
#   my_dl_subsamples,
#   sc,
#   verbose = TRUE
# )
#
# coclustering_results <- calculate_coclustering(
#   batch_subsample_results,
#   sol_df,
#   verbose = TRUE
# )
#
# cocluster_dfs <- coclustering_results$"cocluster_dfs"
#
# cocluster_heatmap(
#   cocluster_dfs[[1]],
#   dl = my_dl,
#   top_hm = list(
#     "Income" = "household_income",
#     "Pubertal Status" = "pubertal_status"
#   ),
#   annotation_colours = list(
#     "Pubertal Status" = colour_scale(
#       c(1, 4),
#       min_colour = "black",
#       max_colour = "purple"
#     ),
#     "Income" = colour_scale(
#       c(0, 4),
#       min_colour = "black",
#       max_colour = "red"
#     )
#   )
# )
# )
# )

```

**Description**

**[Deprecated]** Defunct function for converting a data list into a data frame. Please use `as.data.frame()` instead.

**Usage**

```
collapse_dl(data_list)
```

**Arguments**

`data_list` A nested list of input data from `generate_data_list()`.

**Value**

A "data.frame"-formatted version of the provided data list.

---

colour_scale	<i>Return a colour ramp for a given vector</i>
--------------	--

---

**Description**

Given a numeric vector and min and max colour values, return a colour ramp that assigns a colour to each element in the vector. This function is a wrapper for `circlize::colorRamp2`.

**Usage**

```
colour_scale(data, min_colour, max_colour)
```

**Arguments**

`data` Vector of numeric values.  
`min_colour` Minimum colour value.  
`max_colour` Maximum colour value.

**Value**

A "function" class object that can build a circlize-style colour ramp.

---

config_heatmap	<i>Heatmap for visualizing an SNF config</i>
----------------	--

---

### Description

Create a heatmap where each row corresponds to a different set of hyperparameters in an SNF config object. Numeric parameters are scaled normalized and non-numeric parameters are added as heatmap annotations. Rows can be reordered to match prior meta clustering results.

### Usage

```
config_heatmap(
  sc,
  order = NULL,
  hide_fixed = FALSE,
  show_column_names = TRUE,
  show_row_names = TRUE,
  rect_gp = grid::gpar(col = "black"),
  colour_breaks = c(0, 1),
  colours = c("black", "darkseagreen"),
  column_split_vector = NULL,
  row_split_vector = NULL,
  column_split = NULL,
  row_split = NULL,
  column_title = NULL,
  include_weights = TRUE,
  include_settings = TRUE,
  ...
)
```

### Arguments

sc	An snf_config class object.
order	Numeric vector indicating row ordering of SNF config.
hide_fixed	Whether fixed parameters should be removed.
show_column_names	Whether show column names.
show_row_names	Whether show row names.
rect_gp	Graphic parameters for drawing rectangles (for heatmap body). The value should be specified by <code>gpar</code> and fill parameter is ignored.
colour_breaks	Numeric vector of breaks for the legend.
colours	Vector of colours to use for the heatmap. Should match the length of colour_breaks.
column_split_vector	Vector of indices to split columns by.

```

row_split_vector      Vector of indices to split rows by.
column_split         Split on columns. For heatmap splitting, please refer to https://jokergoo.github.io/ComplexHeatmap-reference/book/a-single-heatmap.html#heatmap-split.
.
row_split            Same as split.
column_title         Title on the column.
include_weights      If TRUE, includes feature weights of the weights matrix into the config heatmap.
include_settings     If TRUE, includes columns from the settings data frame into the config heatmap.
...                  Additional parameters passed to ComplexHeatmap::Heatmap.

```

**Value**

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the scaled values of the provided SNF config.

**Examples**

```

dl <- data_list(
  list(income, "household_income", "demographics", "ordinal"),
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  list(fav_colour, "favourite_colour", "demographics", "categorical"),
  list(anxiety, "anxiety", "behaviour", "ordinal"),
  list(depress, "depressed", "behaviour", "ordinal"),
  uid = "unique_id"
)

sc <- snf_config(
  dl,
  n_solutions = 10,
  dropout_dist = "uniform"
)

config_heatmap(sc)

```

---

cort\_sa

*Mock ABCD cortical surface area data*


---

**Description**

Like the mock data frame "abcd\_cort\_sa", but with "unique\_id" as the "uid".

**Usage**

```
cort_sa
```

**Format**

cort\_sa:

A data frame with 188 rows and 152 columns:

**unique\_id** The unique identifier of the ABCD dataset  
 ... Cortical surface areas of various ROIs (mm<sup>2</sup>, I think)

**Source**

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

cort\_t

*Mock ABCD cortical thickness data*

---

**Description**

Like the mock data frame "abcd\_cort\_t", but with "unique\_id" as the "uid".

**Usage**

cort\_t

**Format**

cort\_t:

A data frame with 188 rows and 152 columns:

**unique\_id** The unique identifier of the ABCD dataset  
 ... Cortical thicknesses of various ROIs (mm<sup>3</sup>, I think)



## Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

data\_list

*Build a data\_list class object*

---

## Description

`data_list()` constructs a data list object which inherits from classes `data_list` and `list`. This object is the primary way in which features to be used along the `metasnf` clustering pipeline are stored. The data list is fundamentally a 2-level nested list object where each inner list contains a data frame and associated metadata for that data frame. The metadata includes the name of the data frame, the 'domain' of that data frame (the broader source of information that the input data frame is capturing, determined by user's domain knowledge), and the type of feature stored in the data frame (continuous, discrete, ordinal, categorical, or mixed).

## Usage

```
data_list(..., uid)
```

## Arguments

... Any number of lists formatted as (df, "df\_name", "df\_domain", "df\_type") and/or any number of lists of lists formatted as (df, "df\_name", "df\_domain", "df\_type").

uid (character) the name of the uid column currently used data. data frame.

## Examples

```
heart_rate_df <- data.frame(
  patient_id = c("1", "2", "3"),
  var1 = c(0.04, 0.1, 0.3),
```

```
    var2 = c(30, 2, 0.3)
  )

  personality_test_df <- data.frame(
    patient_id = c("1", "2", "3"),
    var3 = c(900, 1990, 373),
    var4 = c(509, 2209, 83)
  )

  survey_response_df <- data.frame(
    patient_id = c("1", "2", "3"),
    var5 = c(1, 3, 3),
    var6 = c(2, 3, 3)
  )

  city_df <- data.frame(
    patient_id = c("1", "2", "3"),
    var7 = c("toronto", "montreal", "vancouver")
  )

  # Explicitly (Name each nested list element):
  dl <- data_list(
    list(
      data = heart_rate_df,
      name = "heart_rate",
      domain = "clinical",
      type = "continuous"
    ),
    list(
      data = personality_test_df,
      name = "personality_test",
      domain = "surveys",
      type = "continuous"
    ),
    list(
      data = survey_response_df,
      name = "survey_response",
      domain = "surveys",
      type = "ordinal"
    ),
    list(
      data = city_df,
      name = "city",
      domain = "location",
      type = "categorical"
    ),
    uid = "patient_id"
  )

  # Compact loading
  dl <- data_list(
    list(heart_rate_df, "heart_rate", "clinical", "continuous"),
    list(personality_test_df, "personality_test", "surveys", "continuous"),
```

```

    list(survey_response_df, "survey_response", "surveys", "ordinal"),
    list(city_df, "city", "location", "categorical"),
    uid = "patient_id"
  )

# Printing data list summaries
summary(dl)

# Alternative loading: providing a single list of lists
list_of_lists <- list(
  list(heart_rate_df, "data1", "domain1", "continuous"),
  list(personality_test_df, "data2", "domain2", "continuous")
)

dl <- data_list(
  list_of_lists,
  uid = "patient_id"
)

```

---

depress

*Mock ABCD depression data*


---

### Description

Like the mock data frame "abcd\_depress", but with "unique\_id" as the "uid".

### Usage

```
depress
```

### Format

depress:

A data frame with 275 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**cbcl\_depress\_r** Ordinal value of impairment on CBCL anxiety, either 0 (no impairment), 1 (borderline clinical), or 2 (clinically impaired)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987,

U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

diagnosis_df	<i>Mock diagnosis data</i>
--------------	----------------------------

---

### Description

This is the same data as cancer\_diagnosis\_df, with renamed features and columns.

### Usage

```
diagnosis_df
```

### Format

diagnosis\_df:

A data frame with 200 rows and 2 columns:

**patient\_id** Random three-digit number uniquely identifying the patient

**diagnosis** Mock diagnosis feature

### Source

This data came from the SNFtool package, with slight modifications.

---

dist_fns	<i>Built-in distance functions</i>
----------	------------------------------------

---

### Description

These functions can be used when building a metasnf distance functions list. Each function converts a data frame into to a distance matrix.

**Usage**

```
euclidean_distance(df, weights_row)
gower_distance(df, weights_row)
sn_euclidean_distance(df, weights_row)
sew_euclidean_distance(df, weights_row)
hamming_distance(df, weights_row)
```

**Arguments**

df	Data frame containing at least 1 data column
weights_row	Single-row data frame where the column names contain the column names in df and the row contains the corresponding weights_row.

**Details**

Functions that work for numeric data:

- `euclidean_distance`: typical Euclidean distance
- `sn_euclidean_distance`: Data frame is first standardized and normalized before typical Euclidean distance is applied
- `siw_euclidean_distance`: Squared (including weights) Euclidean distance, where the weights are also squared
- `sew_euclidean_distance`: Squared (excluding weights) Euclidean distance, where the weights are not also squared

Functions that work for binary data:

- `hamming_distance`: typical Hamming distance

Functions that work for any type of data:

- `gower_distance`: Gower distance (`cluster::daisy`)

**Value**

A matrix class object containing pairwise distances.

---

<code>dist_fns_list</code>	<i>Build a distance metrics list</i>
----------------------------	--------------------------------------

---

### Description

The distance metrics list object (inherits classes `dist_fns_list` and `list`) is a list that stores R functions which can convert a data frame of features into a matrix of pairwise distances. The list is a nested one, where the first layer of the list can hold up to 5 items (one for each of the `metasnf` recognized feature types, continuous, discrete, ordinal, categorical, and mixed), and the second layer can hold an arbitrary number of distance functions for each of those types.

### Usage

```
dist_fns_list(
  cnt_dist_fns = NULL,
  dsc_dist_fns = NULL,
  ord_dist_fns = NULL,
  cat_dist_fns = NULL,
  mix_dist_fns = NULL,
  automatic_standard_normalize = FALSE,
  use_default_dist_fns = FALSE
)
```

### Arguments

<code>cnt_dist_fns</code>	A named list of continuous distance metric functions.
<code>dsc_dist_fns</code>	A named list of discrete distance metric functions.
<code>ord_dist_fns</code>	A named list of ordinal distance metric functions.
<code>cat_dist_fns</code>	A named list of categorical distance metric functions.
<code>mix_dist_fns</code>	A named list of mixed distance metric functions.
<code>automatic_standard_normalize</code>	If TRUE, will automatically use standard normalization prior to calculation of any numeric distances. This parameter overrides all other distance functions list-related parameters.
<code>use_default_dist_fns</code>	If TRUE, prepend the base distance metrics (euclidean distance for continuous, discrete, and ordinal data and gower distance for categorical and mixed data) to the resulting distance metrics list.

### Details

Call `?distance_metrics` to see all distance metric functions provided in `metasnf`.

### Value

A distance metrics list object.

**Examples**

```

# Using just the base distance metrics -----
dist_fns_list <- dist_fns_list()

# Adding your own metrics -----
# This will contain only the and user-provided distance function:
cubed_euclidean <- function(df, weights_row) {
  # (your code that converts a data frame to a distance metric here...)
  weights <- diag(weights_row, nrow = length(weights_row))
  weighted_df <- as.matrix(df) %*% weights
  distance_matrix <- weighted_df |>
    stats::dist(method = "euclidean") |>
    as.matrix()
  distance_matrix <- distance_matrix^3
  return(distance_matrix)
}

dist_fns_list <- dist_fns_list(
  cnt_dist_fns = list(
    "my_cubed_euclidean" = cubed_euclidean
  )
)

# Using default base metrics-----
# Call ?distance_metrics to see all distance metric functions provided in
# metasnf. The code below will contain a mix of user-provided and built-in
# distance metric functions.
dist_fns_list <- dist_fns_list(
  cnt_dist_fns = list(
    "my_distance_metric" = cubed_euclidean
  ),
  dsc_dist_fns = list(
    "my_distance_metric" = cubed_euclidean
  ),
  ord_dist_fns = list(
    "my_distance_metric" = cubed_euclidean
  ),
  cat_dist_fns = list(
    "my_distance_metric" = gower_distance
  ),
  mix_dist_fns = list(
    "my_distance_metric" = gower_distance
  ),
  use_default_dist_fns = TRUE
)

```

**Description**

This function enables manipulating a `data_list` class object with `lapply` syntax without removing that object's `data_list` class attribute. The function will only preserve this attribute if the result of the `apply` call has a valid data list structure.

**Usage**

```
dlapply(X, FUN, ...)
```

**Arguments**

X	A <code>data_list</code> class object.
FUN	The function to be applied to each data list component.
...	Optional arguments to FUN.

**Value**

If FUN applied to each component of X yields a valid data list, a data list. Otherwise, a list.

**Examples**

```
# Convert all UID values to lowercase
dl <- data_list(
  list(abcd_income, "income", "demographics", "discrete"),
  list(abcd_colour, "colour", "likes", "categorical"),
  uid = "patient"
)

dl_lower <- dlapply(
  dl,
  function(x) {
    x$"data$"uid <- tolower(x$"data$"uid)
    return(x)
  }
)
```

---

dl\_variable\_summary    *Variable-level summary of a data list*

---

**Description**

**[Deprecated]** Defunct function to summarize a data list. Please use `summary()` with argument `scope = "feature"` instead.

**Usage**

```
dl_variable_summary(dl)
```



**Arguments**

`dl` A nested list of input data from `data_list()`.

**Value**

`variable_level_summary` A data frame containing the name, type, and domain of every variable in a data list.

---

`esm_manhattan_plot` *Manhattan plot of feature-cluster association p-values*

---

**Description**

Manhattan plot of feature-cluster association p-values

**Usage**

```
esm_manhattan_plot(
  esm,
  neg_log_pval_thresh = 5,
  threshold = NULL,
  point_size = 5,
  jitter_width = 0.1,
  jitter_height = 0.1,
  text_size = 15,
  plot_title = NULL,
  hide_x_labels = FALSE,
  bonferroni_line = FALSE
)
```

**Arguments**

`esm` Extended solutions data frame storing associations between features and cluster assignments. See `?extend_solutions`.

`neg_log_pval_thresh` Threshold for negative log p-values.

`threshold` P-value threshold to plot dashed line at.

`point_size` Size of points in the plot.

`jitter_width` Width of jitter.

`jitter_height` Height of jitter.

`text_size` Size of text in the plot.

`plot_title` Title of the plot.

`hide_x_labels` If TRUE, hides x-axis labels.

`bonferroni_line` If TRUE, plots a dashed black line at the Bonferroni-corrected equivalent of the p-value threshold.

**Value**

A Manhattan plot (class "gg", "ggplot") showing the association p-values of features against each solution in the provided solutions data frame.

**Examples**

```
# full_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   list(anxiety, "anxiety", "behaviour", "ordinal"),
#   list(depress, "depressed", "behaviour", "ordinal"),
#   uid = "unique_id"
# )
#
# dl <- full_dl[1:3]
# target_dl <- full_dl[4:5]
#
# set.seed(42)
# sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
# )
#
# sol_df <- batch_snf(dl, sc)
#
# ext_sol_df <- extend_solutions(
#   sol_df,
#   dl = dl,
#   target = target_dl,
#   min_pval = 1e-10 # p-values below 1e-10 will be thresholded to 1e-10
# )
#
# esm_manhattan <- esm_manhattan_plot(
#   ext_sol_df[1:5, ],
#   neg_log_pval_thresh = 5,
#   threshold = 0.05,
#   point_size = 3,
#   jitter_width = 0.1,
#   jitter_height = 0.1,
#   plot_title = "Feature-Solution Associations",
#   text_size = 14,
#   bonferroni_line = TRUE
# )
```

---

estimate\_nclust\_given\_graph

*Estimate number of clusters for a similarity matrix*

---

**Description**

Calculate eigengap and rotation-cost estimates of the number of clusters to use when clustering a similarity matrix. This function was adapted from `SNFtool::estimateClustersGivenGraph`, but scales up the Laplacian operator prior to eigenvalue calculations to minimize the risk of floating point-related errors.

**Usage**

```
estimate_nclust_given_graph(W, NUMC = 2:10)
```

**Arguments**

W	Similarity matrix to calculate number of clusters for.
NUMC	Range of cluster counts to consider among when picking best number of clusters.

**Value**

A list containing the top two eigengap and rotation-cost estimates for the number of clusters in a given similarity matrix.

**Examples**

```
input_dl <- data_list(
  list(gender_df, "gender", "demographics", "categorical"),
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),
  uid = "patient_id"
)

sc <- snf_config(input_dl, n_solutions = 1)
sol_df <- batch_snf(input_dl, sc, return_sim_mats = TRUE)
sim_mat <- sim_mats_list(sol_df)[[1]]
estimate_nclust_given_graph(sim_mat)
```

---

expression\_df

*Modification of SNFtool mock data frame "Data1"*

---

**Description**

Modification of SNFtool mock data frame "Data1"

**Usage**

```
expression_df
```

**Format**

expression\_df:

A data frame with 200 rows and 3 columns:

**gene\_1\_expression** Mock gene expression feature

**gene\_2\_expression** Mock gene expression feature

**patient\_id** Random three-digit number uniquely identifying the patient

**Source**

This data came from the SNFtool package, with slight modifications.

---

extend_solutions	<i>Extend a solutions data frame to include outcome evaluations</i>
------------------	---

---

**Description**

Extend a solutions data frame to include outcome evaluations

**Usage**

```
extend_solutions(
  sol_df,
  target_dl = NULL,
  dl = NULL,
  cat_test = "chi_squared",
  min_pval = 1e-10,
  processes = 1,
  verbose = FALSE
)
```

**Arguments**

sol_df	Result of batch_snf storing cluster solutions and the settings that were used to generate them.
target_dl	A data list with features to calculate p-values for. Features in the target list will be included during p-value summary measure calculations.
dl	A data list with features to calculate p-values for, but that should not be incorporated into p-value summary measure columns (i.e., min/mean/max p-value columns).
cat_test	String indicating which statistical test will be used to associate cluster with a categorical feature. Options are "chi_squared" for the Chi-squared test and "fisher_exact" for Fisher's exact test.
min_pval	If assigned a value, any p-value less than this will be replaced with this value.
processes	The number of processes to use for parallelization. Progress is only reported for sequential processing (processes = 1).
verbose	If TRUE, output progress to console.

**Value**

An extended solutions data frame (`ext_sol_df` class object) that contains p-value columns for each outcome in the provided data lists

**Examples**

```
input_dl <- data_list(
  list(gender_df, "gender", "demographics", "categorical"),
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),
  uid = "patient_id"
)

sc <- snf_config(input_dl, n_solutions = 2)

sol_df <- batch_snf(input_dl, sc)

ext_sol_df <- extend_solutions(sol_df, input_dl)
```

---

fav_colour	<i>Mock ABCD "colour" data</i>
------------	--------------------------------

---

**Description**

Like the mock data frame "abcd\_colour", but with "unique\_id" as the "uid".

**Usage**

```
fav_colour
```

**Format**

**fav\_colour:**

A data frame with 275 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**colour** Categorical transformation of `cbcl_depress`.

**Source**

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988,

U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

features	<i>Return character vector of features stored in an object</i>
----------	--

---

### Description

Return character vector of features stored in an object

### Usage

```
features(x)
```

### Arguments

x                    The object to pull features from.

### Value

A character vector of features in x.

---

gender_df	<i>Mock gender data</i>
-----------	-------------------------

---

### Description

Mock gender data

### Usage

```
gender_df
```

### Format

gender\_df:

A data frame with 200 rows and 2 columns:

**patient\_id** Random three-digit number uniquely identifying the patient

**gender\_df** Mock gene methylation feature

### Source

This data came from the SNFtool package, with slight modifications.

---

```
generate_clust_algs_list
```

*Generate a clustering algorithms list*

---

### Description

**[Deprecated]** Deprecated function for building a clustering algorithms list. Please use `clust_fns_list()` (or better yet, `snf_config()`) instead.

### Usage

```
generate_clust_algs_list(..., disable_base = FALSE)
```

### Arguments

<code>...</code>	An arbitrary number of named clustering functions
<code>disable_base</code>	If TRUE, do not prepend the base clustering algorithms ( <code>spectral_eigen</code> and <code>spectral_rot</code> , which apply spectral clustering and use the eigen-gap and rotation cost heuristics respectively for determining the number of clusters in the graph.

### Value

A list of clustering algorithm functions that can be passed into the `batch_snf` and `generate_settings_list` functions.

---

```
generate_distance_metrics_list
```

*Generate a list of distance metrics*

---

### Description

**[Deprecated]** Deprecated function for building a distance metrics list. Please use `dist_fns_list()` (or better yet, `snf_config()`) instead.

### Usage

```
generate_distance_metrics_list(
  continuous_distances = NULL,
  discrete_distances = NULL,
  ordinal_distances = NULL,
  categorical_distances = NULL,
  mixed_distances = NULL,
  keep_defaults = TRUE
)
```

**Arguments**

continuous_distances	A named list of distance metric functions
discrete_distances	A named list of distance metric functions
ordinal_distances	A named list of distance metric functions
categorical_distances	A named list of distance metric functions
mixed_distances	A named list of distance metric functions
keep_defaults	If TRUE (default), prepend the base distance metrics (euclidean and standard normalized euclidean)

**Value**

A nested and named list of distance metrics functions.

---

generate\_settings\_matrix

*Build a settings data frame*

---

**Description**

**[Deprecated]** Deprecated function for building a settings matrix. Please use settings\_df() instead.

**Usage**

```
generate_settings_matrix(...)
```

**Arguments**

... Arguments used to generate a settings matrix.

**Value**

Raises a deprecated error.



---

get_clusters	<i>Extract cluster membership vector from one solutions data frame row</i>
--------------	--

---

### Description

**[Deprecated]** Deprecated function for building extracting cluster solutions from a solutions data frame. Please use `t()` instead.

This function takes in a single row of a solutions data frame and returns a vector containing the cluster assignments for each observation. It is similar to `get_cluster_df()`, which takes a solutions data frame with only one row and returns a data frame with two columns: "cluster" and "uid" (the UID of the observation) and `get_cluster_solutions()`, which takes a solutions data frame with any number of rows and returns a data frame indicating the cluster assignments for each of those rows.

### Usage

```
get_clusters(sol_df_row)
```

### Arguments

`sol_df_row`      Output matrix row.

### Value

clusters Vector of assigned clusters.

---

get_cluster_df	<i>Extract cluster membership information from one solutions data frame row</i>
----------------	---

---

### Description

**[Deprecated]** Deprecated function for building extracting cluster solutions from a solutions data frame. Please use `t()` instead.

This function takes in a single row of a solutions data frame and returns a data frame containing the cluster assignments for each uid. It is similar to `get_clusters()`, which takes one solutions data frame row and returns a vector of cluster assignments' and `get_cluster_solutions()`, which takes a solutions data frame with any number of rows and returns a data frame indicating the cluster assignments for each of those rows.

### Usage

```
get_cluster_df(sol_df_row)
```

**Arguments**

sol\_df\_row      One row from a solutions data frame.

**Value**

cluster\_df data frame of cluster and uid.

---

get\_cluster\_solutions    *Extract cluster membership information from a sol\_df*

---

**Description**

**[Deprecated]** Deprecated function for building extracting cluster solutions from a solutions data frame. Please use `t()` instead.

This function takes in a solutions data frame and returns a data frame containing the cluster assignments for each uid. It is similar to `'get_clusters()'`, which takes one solutions data frame row and returns a vector of cluster assignments' and `get_cluster_df()`, which takes a solutions matrix with only one row and returns a data frame with two columns: "cluster" and "uid" (the UID of the observation).

**Usage**

```
get_cluster_solutions(sol_df)
```

**Arguments**

sol\_df            A sol\_df.

**Value**

A "data.frame" object where each row is an observation and each column (apart from the uid column) indicates the cluster that observation was assigned to for the corresponding solutions data frame row.

---

get\_complete\_uids      *Pull complete-data UIDs from a list of data frames*

---

**Description**

This function identifies all observations within a list of data frames that have no missing data across all data frames. This function is useful when constructing data lists of distinct feature sets from the same sample of observations. As `data_list()` strips away observations with any missing data, distinct sets of observations may be generated by building a data list from the same group of observations over different sets of features. Reducing the pool of observations to only those with complete UIDs first will avoid downstream generation of data lists of differing sizes.

**Usage**

```
get_complete_uids(list_of_dfs, uid)
```

**Arguments**

```
list_of_dfs    List of data frames.  
uid            Name of column across data frames containing UIDs
```

**Value**

A character vector of the UIDs of observations that have complete data across the provided list of data frames.

**Examples**

```
complete_uids <- get_complete_uids(  
  list(income, pubertal, anxiety, depress),  
  uid = "unique_id"  
)  
  
income <- income[income$"unique_id" %in% complete_uids, ]  
pubertal <- pubertal[pubertal$"unique_id" %in% complete_uids, ]  
anxiety <- anxiety[anxiety$"unique_id" %in% complete_uids, ]  
depress <- depress[depress$"unique_id" %in% complete_uids, ]  
  
input_dl <- data_list(  
  list(income, "income", "demographics", "ordinal"),  
  list(pubertal, "pubertal", "demographics", "continuous"),  
  uid = "unique_id"  
)  
  
target_dl <- data_list(  
  list(anxiety, "anxiety", "behaviour", "ordinal"),  
  list(depress, "depressed", "behaviour", "ordinal"),  
  uid = "unique_id"  
)
```

---

get\_dl\_uids

*Extract UIDs from a data list*

---

**Description**

**[Deprecated]** Deprecated function for extracting UIDs from a data list. Please use `uids()` instead.

**Usage**

```
get_dl_uids(dl, prefix = FALSE)
```

**Arguments**

- dl                    A nested list of input data from `data_list()`.
- prefix                If TRUE, preserves the "uid\_" prefix added to UIDs when creating a data list.

**Value**

A character vector of the UID labels contained in a data list.

---

get\_heatmap\_order     *Return the row or column ordering present in a heatmap*

---

**Description**

Return the row or column ordering present in a heatmap

**Usage**

```
get_heatmap_order(heatmap, type = "rows")
```

**Arguments**

- heatmap              A heatmap object to collect ordering from.
- type                  The type of ordering to return. Either "rows" or "columns".

**Value**

A numeric vector of the ordering used within the provided ComplexHeatmap "Heatmap" object.

---

get\_matrix\_order     *Return the hierarchical clustering order of a matrix*

---

**Description**

Return the hierarchical clustering order of a matrix

**Usage**

```
get_matrix_order(matrix, dist_method = "euclidean", hclust_method = "complete")
```

**Arguments**

matrix	Matrix to cluster.
dist_method	Distance method to use when calculating sorting order to of the matrix. Argument is directly passed into stats::dist. Options include "euclidean", "maximum", "manhattan", "canberra", "binary", or "minkowski".
hclust_method	Agglomerative method to use when calculating sorting order by stats::hclust. Options include "ward.D", "ward.D2", "single", "complete", "average", "mcquitty", "median", or "centroid".

**Value**

A numeric vector of the ordering derived by the specified hierarchical clustering method applied to the provided matrix.

**Examples**

```
# dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   list(anxiety, "anxiety", "behaviour", "ordinal"),
#   list(depress, "depressed", "behaviour", "ordinal"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
# )
#
# sol_df <- batch_snf(dl, sc)
#
# ext_sol_df <- extend_solutions(
#   sol_df,
#   dl = dl,
#   min_pval = 1e-10 # p-values below 1e-10 will be thresholded to 1e-10
# )
#
# # Calculate pairwise similarities between cluster solutions
# sol_aris <- calc_aris(sol_df)
#
# # Extract hierarchical clustering order of the cluster solutions
# meta_cluster_order <- get_matrix_order(sol_aris)
```

---

 get\_pvals

*Get p-values from an extended solutions data frame*


---

### Description

This function can be used to neatly format the p-values associated with an extended solutions data frame. It can also calculate the negative logs of those p-values to make it easier to interpret large-scale differences.

### Usage

```
get_pvals(ext_sol_df, negative_log = FALSE, keep_summaries = TRUE)
```

### Arguments

`ext_sol_df` The output of `extend_solutions`. A data frame that contains at least one p-value column ending in "\_pval".

`negative_log` If TRUE, will replace p-values with negative log p-values.

`keep_summaries` If FALSE, will remove the mean, min, and max p-value.

### Value

A "data.frame" class object Of only the p-value related columns of the provided `ext_sol_df`.

---

 get\_representative\_solutions

*Extract representative solutions from a matrix of ARIs*


---

### Description

Following clustering with `batch_snf`, a matrix of pairwise ARIs that show how related each cluster solution is to each other can be generated by the `calc_aris` function. Partitioning of the ARI matrix can be done by visual inspection of `meta_cluster_heatmap()` results or by `shiny_annotator`. Given the indices of meta cluster boundaries, this function will return a single representative solution from each meta cluster based on maximum average ARI to all other solutions within that meta cluster.

### Usage

```
get_representative_solutions(aris, sol_df, filter_fn = NULL)
```

**Arguments**

aris	Matrix of adjusted rand indices from calc_aris()
sol_df	Output of batch_snf containing cluster solutions.
filter_fn	Optional function to filter the meta-cluster by prior to maximum average ARI determination. This can be useful if you are explicitly trying to select a solution that meets a certain condition, such as only picking from the 4 cluster solutions within a meta cluster. An example valid function could be <code>fn &lt;- function(x) x[<math>x</math>["nclust"] == 4, ]</code> .

**Value**

The provided solutions data frame reduced to just one row per meta cluster defined by the split vector.

**Examples**

```
# dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   list(anxiety, "anxiety", "behaviour", "ordinal"),
#   list(depress, "depressed", "behaviour", "ordinal"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
# )
#
# sol_df <- batch_snf(dl, sc)
#
# ext_sol_df <- extend_solutions(
#   sol_df,
#   dl = dl,
#   min_pval = 1e-10 # p-values below 1e-10 will be thresholded to 1e-10
# )
#
# # Calculate pairwise similarities between cluster solutions
# sol_aris <- calc_aris(sol_df)
#
# # Extract hierarchical clustering order of the cluster solutions
# meta_cluster_order <- get_matrix_order(sol_aris)
#
# # Identify meta cluster boundaries with shiny app or trial and error
# # ari_hm <- meta_cluster_heatmap(sol_aris, order = meta_cluster_order)
# # shiny_annotator(ari_hm)
#
# # Result of meta cluster examination
```

```
# split_vec <- c(2, 5, 12, 17)
#
# ext_sol_df <- label_meta_clusters(ext_sol_df, split_vec, meta_cluster_order)
#
# # Extracting representative solutions from each defined meta cluster
# rep_solutions <- get_representative_solutions(sol_aris, ext_sol_df)
```

---

income

*Mock ABCD income data*

---

### Description

Like the mock data frame "abcd\_h\_income", but with "unique\_id" as the "uid".

Like the mock data frame "abcd\_cort\_sa", but with "unique\_id" as the "uid".

### Usage

```
income
```

```
income
```

### Format

income:

A data frame with 300 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**household\_income** Household income in 3 category levels (low = 1, medium = 2, high = 3)

income:

A data frame with 300 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**household\_income** Household income in 3 category levels (low = 1, medium = 2, high = 3)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148,



U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

is\_data\_list

*Test if the object is a data list*

---

### **Description**

Given an object, returns TRUE if that object inherits from the data\_list class.

### **Usage**

```
is_data_list(x)
```

### **Arguments**

x                    An object.

### **Value**

TRUE if the object inherits from the data\_list class.

---

jitter_plot	<i>Jitter plot separating a feature by cluster</i>
-------------	--

---

**Description**

Jitter plot separating a feature by cluster

**Usage**

```
jitter_plot(df, feature)
```

**Arguments**

df	A data.frame containing cluster column and the feature to plot.
feature	The feature to plot.

**Value**

A jitter+violin plot (class "gg", "ggplot") showing the distribution of a feature across clusters.

---

label_meta_clusters	<i>Assign meta cluster labels to rows of a solutions data frame or extended solutions data frame</i>
---------------------	--

---

**Description**

Given a solutions data frame or extended solutions data frame class object and a numeric vector indicating which rows correspond to which meta clusters, assigns meta clustering information to the "meta\_clusters" attribute of the data frame.

**Usage**

```
label_meta_clusters(sol_df, split_vector, order = NULL)
```

**Arguments**

sol_df	A solutions data frame or extended solutions data frame to assign meta clusters to.
split_vector	A numeric vector indicating which rows of sol_df should be the split points for meta cluster labeling.
order	An optional numeric vector indicating how the solutions data frame should be reordered prior to meta cluster labeling. This vector can be obtained by running <code>get_matrix_order()</code> on an ARI matrix, which itself can be obtained by calling <code>calc_aris()</code> on a solutions data frame.

**Value**

A solutions data frame with a populated "meta\_clusters" attribute.

**Examples**

```
#dl <- data_list(
#   list(cort_sa, "cortical_surface_area", "neuroimaging", "continuous"),
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
#)
#
#set.seed(42)
#my_sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
#)
#
#sol_df <- batch_snf(dl, my_sc)
#
#sol_df
#
#sol_aris <- calc_aris(sol_df)
#
#meta_cluster_order <- get_matrix_order(sol_aris)
#
## `split_vec` found by iteratively plotting ari_hm or by ?shiny_annotator()
#split_vec <- c(6, 10, 16)
#ari_hm <- meta_cluster_heatmap(
#   sol_aris,
#   order = meta_cluster_order,
#   split_vector = split_vec
#)
#
#mc_sol_df <- label_meta_clusters(
#   sol_df,
#   order = meta_cluster_order,
#   split_vector = split_vec
#)
#
#mc_sol_df
```

**Description**

Given a solutions data frame containing clustered observations and a data list containing those clustered observations as well as additional to-be-clustered observations, this function will re-run SNF to generate a similarity matrix of all observations and use the label propagation algorithm to assigned predicted clusters to the unclustered observations.

**Usage**

```
label_propagate(partial_sol_df, full_dl, verbose = FALSE)
```

**Arguments**

`partial_sol_df` A solutions data frame derived from the training set.  
`full_dl` A data list containing observations from both the training and testing sets.  
`verbose` If TRUE, output progress to console.

**Value**

A data frame with one row per observation containing a column for UIDs, a column for whether the observation was in the train (original) or test (held out) set, and one column per row of the solutions data frame indicating the original and propagated clusters.

**Examples**

```
## Function to identify observations with complete data
#uids_with_complete_obs <- get_complete_uids(
#  list(subc_v, income, pubertal, anxiety, depress),
#  uid = "unique_id"
#)
#
## Dataframe assigning 80% of observations to train and 20% to test
#train_test_split <- train_test_assign(
#  train_frac = 0.8,
#  uids = uids_with_complete_obs
#)
#
## Pulling the training and testing observations specifically
#train_obs <- train_test_split$"train"
#test_obs <- train_test_split$"test"
#
## Partition a training set
#train_subc_v <- subc_v[subc_v$"unique_id" %in% train_obs, ]
#train_income <- income[income$"unique_id" %in% train_obs, ]
#train_pubertal <- pubertal[pubertal$"unique_id" %in% train_obs, ]
#train_anxiety <- anxiety[anxiety$"unique_id" %in% train_obs, ]
#train_depress <- depress[depress$"unique_id" %in% train_obs, ]
#
## Partition a test set
#test_subc_v <- subc_v[subc_v$"unique_id" %in% test_obs, ]
#test_income <- income[income$"unique_id" %in% test_obs, ]
```

```

#test_pubertal <- pubertal[pubertal$"unique_id" %in% test_obs, ]
#test_anxiety <- anxiety[anxiety$"unique_id" %in% test_obs, ]
#test_depress <- depress[depress$"unique_id" %in% test_obs, ]
#
## Find cluster solutions in the training set
#train_dl <- data_list(
#  list(train_subc_v, "subc_v", "neuroimaging", "continuous"),
#  list(train_income, "household_income", "demographics", "continuous"),
#  list(train_pubertal, "pubertal_status", "demographics", "continuous"),
#  uid = "unique_id"
#)
#
## We'll pick a solution that has good separation over our target features
#train_target_dl <- data_list(
#  list(train_anxiety, "anxiety", "behaviour", "ordinal"),
#  list(train_depress, "depressed", "behaviour", "ordinal"),
#  uid = "unique_id"
#)
#
#sc <- snf_config(
#  train_dl,
#  n_solutions = 5,
#  min_k = 10,
#  max_k = 30
#)
#
#train_sol_df <- batch_snf(
#  train_dl,
#  sc,
#  return_sim_mats = TRUE
#)
#
#ext_sol_df <- extend_solutions(
#  train_sol_df,
#  train_target_dl
#)
#
## Determining solution with the lowest minimum p-value
#lowest_min_pval <- min(ext_sol_df$"min_pval")
#which(ext_sol_df$"min_pval" == lowest_min_pval)
#top_row <- ext_sol_df[1, ]
#
## Propagate that solution to the observations in the test set
## data list below has both training and testing observations
#full_dl <- data_list(
#  list(subc_v, "subc_v", "neuroimaging", "continuous"),
#  list(income, "household_income", "demographics", "continuous"),
#  list(pubertal, "pubertal_status", "demographics", "continuous"),
#  uid = "unique_id"
#)
#
## Use the solutions data frame from the training observations and the data list
## from the training and testing observations to propagate labels to the test observations

```

```
#propagated_labels <- label_propagate(top_row, full_dl)
#
#propagated_labels_all <- label_propagate(ext_sol_df, full_dl)
#
#head(propagated_labels_all)
#tail(propagated_labels_all)
```

---

linear\_adjust

*Linearly correct data list by features with unwanted signal*


---

### Description

Given a data list to correct and another data list of categorical features to linearly adjust for, corrects the first data list based on the residuals of the linear model relating the numeric features in the first data list to the unwanted signal features in the second data list.

### Usage

```
linear_adjust(dl, unwanted_signal_list, sig_digs = NULL)
```

### Arguments

**dl** A nested list of input data from `data_list()`.

**unwanted\_signal\_list** A data list of categorical features that should have their mean differences removed in the first data list.

**sig\_digs** Number of significant digits to round the residuals to.

### Value

A data list ("list") in which each data component has been converted to contain residuals off of the linear model built against the features in the `unwanted_signal_list`.

### Examples

```
has_tutor <- sample(c(1, 0), size = 9, replace = TRUE)
math_score <- 70 + 30 * has_tutor + rnorm(9, mean = 0, sd = 5)

math_df <- data.frame(uid = paste0("id_", 1:9), math = math_score)
tutor_df <- data.frame(uid = paste0("id_", 1:9), tutor = has_tutor)

dl <- data_list(
  list(math_df, "math_score", "school", "continuous"),
  uid = "uid"
)

adjustment_dl <- data_list(
  list(tutor_df, "tutoring", "school", "categorical"),
  uid = "uid"
```

```

)

adjusted_dl <- linear_adjust(dl, adjustment_dl)

adjusted_dl[[1]]$"data" $"math"

# Equivalent to:
as.numeric(resid(lm(math_score ~ has_tutor)))

```

---

mc\_manhattan\_plot      *Manhattan plot of feature-meta cluster association p-values*

---

### Description

Given a data frame of representative meta cluster solutions (see `get_representative_solutions()`), returns a Manhattan plot for showing feature separation across all features in provided data/target lists.

### Usage

```

mc_manhattan_plot(
  ext_sol_df,
  dl = NULL,
  target_dl = NULL,
  variable_order = NULL,
  neg_log_pval_thresh = 5,
  threshold = NULL,
  point_size = 5,
  text_size = 20,
  plot_title = NULL,
  xints = NULL,
  hide_x_labels = FALSE,
  domain_colours = NULL
)

```

### Arguments

<code>ext_sol_df</code>	A <code>sol_df</code> that contains " <code>_pval</code> " columns containing the values to be plotted. This object is the output of <code>extend_solutions()</code> .
<code>dl</code>	List of data frames containing data information.
<code>target_dl</code>	List of data frames containing target information.
<code>variable_order</code>	Order of features to be displayed in the plot.
<code>neg_log_pval_thresh</code>	Threshold for negative log p-values.
<code>threshold</code>	p-value threshold to plot horizontal dashed line at.
<code>point_size</code>	Size of points in the plot.

**text\_size**        Size of text in the plot.  
**plot\_title**      Title of the plot.  
**xints**            Either "outcomes" or a vector of numeric values to plot vertical lines at.  
**hide\_x\_labels**    If TRUE, hides x-axis labels.  
**domain\_colours**   Named vector of colours for domains.

### Value

A Manhattan plot (class "gg", "ggplot") showing the association p-values of features against each solution in the provided solutions data frame, stratified by meta cluster label.

### Examples

```

# dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   list(anxiety, "anxiety", "behaviour", "ordinal"),
#   list(depress, "depressed", "behaviour", "ordinal"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
# )
#
# sol_df <- batch_snf(dl, sc)
#
# ext_sol_df <- extend_solutions(
#   sol_df,
#   dl = dl,
#   min_pval = 1e-10 # p-values below 1e-10 will be thresholded to 1e-10
# )
#
# # Calculate pairwise similarities between cluster solutions
# sol_aris <- calc_aris(sol_df)
#
# # Extract hierarchical clustering order of the cluster solutions
# meta_cluster_order <- get_matrix_order(sol_aris)
#
# # Identify meta cluster boundaries with shiny app or trial and error
# # ari_hm <- meta_cluster_heatmap(sol_aris, order = meta_cluster_order)
# # shiny_annotator(ari_hm)
#
# # Result of meta cluster examination
# split_vec <- c(2, 5, 12, 17)
#
# ext_sol_df <- label_meta_clusters(ext_sol_df, split_vec, meta_cluster_order)

```



```
#
# # Extracting representative solutions from each defined meta cluster
# rep_solutions <- get_representative_solutions(sol_aris, ext_sol_df)
#
# mc_manhattan <- mc_manhattan_plot(
#   rep_solutions,
#   dl = dl,
#   point_size = 3,
#   text_size = 12,
#   plot_title = "Feature-Meta Cluster Associations",
#   threshold = 0.05,
#   neg_log_pval_thresh = 5
# )
#
# mc_manhattan
```

---

merge.snf_config	<i>Merge method for SNF config objects</i>
------------------	--

---

## Description

Merge method for SNF config objects

## Usage

```
## S3 method for class 'snf_config'
merge(x, y, reset_indices = TRUE, ...)
```

## Arguments

x	SNF config to merge.
y	SNF config to merge.
reset_indices	If TRUE (default), re-labels the "solutions" indices in the config from 1 to the number of defined settings.
...	Additional arguments passed into merge function.

## Value

An SNF config combining the rows of both prior configurations.

---

merge_df_list	<i>Merge list of data frames into a single data frame</i>
---------------	---

---

**Description**

This helper function combines all data frames in a single-level list into a single data frame.

**Usage**

```
merge_df_list(df_list, join = "inner", uid = "uid", no_na = FALSE)
```

**Arguments**

df_list	list of data frames.
join	String indicating if join should be "inner" or "full".
uid	Column name to join on. Default is "uid".
no_na	Whether to remove NA values from the merged data frame.

**Value**

Inner join of all data frames in list.

**Examples**

```
merge_df_list(list(income, pubertal), uid = "unique_id")
```

---

merge_dls	<i>Horizontally merge compatible data lists</i>
-----------	---

---

**Description**

Join two data lists with the same components (data frames) but separate observations. To instead merge two data lists that have the same observations but different components, simply use `c()`.

**Usage**

```
merge_dls(dl_1, dl_2)
```

**Arguments**

dl_1	The first data list to merge.
dl_2	The second data list to merge.

**Value**

A data list ("list"-class object) containing the observations of both provided data lists.

---

meta\_cluster\_heatmap *Heatmap of pairwise adjusted rand indices between solutions*

---

## Description

Heatmap of pairwise adjusted rand indices between solutions

## Usage

```
meta_cluster_heatmap(
  aris,
  order = NULL,
  cluster_rows = FALSE,
  cluster_columns = FALSE,
  log_graph = FALSE,
  scale_diag = "none",
  min_colour = "#282828",
  max_colour = "firebrick2",
  col = circlize::colorRamp2(c(min(aris), max(aris)), c(min_colour, max_colour)),
  ...
)
```

## Arguments

aris	Matrix of adjusted rand indices from calc_aris()
order	Numeric vector containing row order of the heatmap.
cluster_rows	Whether rows should be clustered.
cluster_columns	Whether columns should be clustered.
log_graph	If TRUE, log transforms the graph.
scale_diag	Method of rescaling matrix diagonals. Can be "none" (don't change diagonals), "mean" (replace diagonals with average value of off-diagonals), or "zero" (replace diagonals with 0).
min_colour	Colour used for the lowest value in the heatmap.
max_colour	Colour used for the highest value in the heatmap.
col	Colour ramp to use for the heatmap.
...	Additional parameters passed to similarity_matrix_heatmap(), the function that this function wraps.

## Value

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the pairwise adjusted Rand indices (similarities) between the cluster solutions of the provided solutions data frame.

**Examples**

```

#dl <- data_list(
#   list(cort_sa, "cortical_surface_area", "neuroimaging", "continuous"),
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
#)
#
#set.seed(42)
#my_sc <- snf_config(
#   dl = dl,
#   n_solutions = 20,
#   min_k = 20,
#   max_k = 50
#)
#
#sol_df <- batch_snf(dl, my_sc)
#
#sol_df
#
#sol_aris <- calc_aris(sol_df)
#
#meta_cluster_order <- get_matrix_order(sol_aris)
#
## `split_vec` found by iteratively plotting ari_hm or by ?shiny_annotator()
#split_vec <- c(6, 10, 16)
#ari_hm <- meta_cluster_heatmap(
#   sol_aris,
#   order = meta_cluster_order,
#   split_vector = split_vec
#)

```

---

methylation\_df

*Modification of SNFtool mock data frame "Data2"*


---

**Description**

Modification of SNFtool mock data frame "Data2"

**Usage**

```
methylation_df
```

**Format**

```
methylation_df:
```

A data frame with 200 rows and 3 columns:

**gene\_1\_expression** Mock gene methylation feature

**gene\_2\_expression** Mock gene methylation feature  
**patient\_id** Random three-digit number uniquely identifying the patient

### Source

This data came from the SNFtool package, with slight modifications.

---

new\_solutions\_df      *Constructor for solutions\_df class object*

---

### Description

Constructor for solutions\_df class object

### Usage

```
new_solutions_df(sol_df1)
```

### Arguments

sol\_df1      A solutions data frame-like object to be validated and converted into a solutions data frame.

### Value

A solutions\_df class object.

---

n\_features      *Extract number of features stored in an object*

---

### Description

Extract number of features stored in an object

### Usage

```
n_features(x)
```

### Arguments

x      The object to extract number of features from.

### Value

The number of features in x.

---

n_observations	<i>Extract number of observations stored in an object</i>
----------------	---

---

**Description**

Extract number of observations stored in an object

**Usage**

```
n_observations(x)
```

**Arguments**

x                    The object to extract number of observations from.

**Value**

The number of observations in x.

---

print.ari_matrix	<i>Print method for class ari_matrix</i>
------------------	--

---

**Description**

Custom formatted print for weights matrices that outputs information about feature weights functions to the console.

**Usage**

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

**Arguments**

x                    A `ari_matrix` class object.  
...                  Other arguments passed to `print` (not used in this function)

**Value**

Function prints to console but does not return any value.

---

print.clust\_fns\_list    *Print method for class clust\_fns\_list*

---

**Description**

Custom formatted print for clustering functions list objects that outputs information about the contained clustering functions to the console.

**Usage**

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

**Arguments**

x	A clust_fns_list class object.
...	Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

print.data\_list        *Print method for class data\_list*

---

**Description**

Custom formatted print for data list objects that outputs information about the contained observations and components to the console.

**Usage**

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

**Arguments**

x	A data_list class object.
...	Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

```
print.dist_fns_list
```

*Print method for class dist\_fns\_list*

---

**Description**

Custom formatted print for distance metrics list objects that outputs information about the contained distance metrics to the console.

**Usage**

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

**Arguments**

x	A dist_fns_list class object.
...	Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

```
print.ext_solutions_df
```

*Print method for class ext\_solutions\_df*

---

**Description**

Custom formatted print for extended solutions data frame class objects.

**Usage**

```
## S3 method for class 'ext_solutions_df'  
print(x, n = NULL, ...)
```

**Arguments**

x	A ext_solutions_df class object.
n	Number of rows to print, passed into <code>tibble::print.tbl_df()</code> .
...	Other arguments passed to print (not used in this function).

**Value**

Function prints to console but does not return any value.



---

print.settings\_df      *Print method for class settings\_df*

---

**Description**

Custom formatted print for settings data frame that outputs information about SNF hyperparameters to the console.

**Usage**

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

**Arguments**

x                    A settings\_df class object.  
...                  Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

print.snf\_config      *Print method for class snf\_config*

---

**Description**

Custom formatted print for SNF config

**Usage**

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

**Arguments**

x                    A snf\_config class object.  
...                  Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

```
print.solutions_df      Print method for class weights_matrix
```

---

**Description**

Custom formatted print for weights matrices that outputs information about feature weights functions to the console.

**Usage**

```
## S3 method for class 'solutions_df'
print(x, n = NULL, tips = TRUE, ...)
```

**Arguments**

x	A <code>weights_matrix</code> class object.
n	Number of rows to print, passed into <code>tibble::print.tbl_df()</code> .
tips	If TRUE, include lines on how to print more rows / transposed.
...	Other arguments passed to <code>print</code> (not used in this function).

**Value**

Function prints to console but does not return any value.

---

```
print.t_ext_solutions_df
      Print method for class t_ext_solutions_df
```

---

**Description**

Custom formatted print for transposed solutions data frame class objects.

**Usage**

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

**Arguments**

x	A <code>t_solutions_df</code> class object.
...	Other arguments passed to <code>print</code> (not used in this function)

**Value**

Function prints to console but does not return any value.

---

print.t\_solutions\_df *Print method for class t\_solutions\_df*

---

**Description**

Custom formatted print for transposed solutions data frame class objects.

**Usage**

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

**Arguments**

x                    A t\_solutions\_df class object.  
...                  Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

print.weights\_matrix *Print method for class weights\_matrix*

---

**Description**

Custom formatted print for weights matrices that outputs information about feature weights functions to the console.

**Usage**

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

**Arguments**

x                    A weights\_matrix class object.  
...                  Other arguments passed to print (not used in this function)

**Value**

Function prints to console but does not return any value.

---

pubertal	<i>Mock ABCD pubertal status data</i>
----------	---------------------------------------

---

### Description

Like the mock data frame "abcd\_pubertal", but with "unique\_id" as the "uid".

### Usage

```
pubertal
```

### Format

```
pubertal:
```

A data frame with 275 rows and 2 columns:

**unique\_id** The unique identifier of the ABCD dataset

**pubertal\_status** Average reported pubertal status between child and parent (1-5 categorical scale)

### Source

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

pval_heatmap	<i>Heatmap of p-values</i>
--------------	----------------------------

---

### Description

Heatmap of p-values

**Usage**

```

pval_heatmap(
  ext_sol_df,
  order = NULL,
  cluster_columns = TRUE,
  cluster_rows = FALSE,
  show_row_names = FALSE,
  show_column_names = TRUE,
  min_colour = "red2",
  max_colour = "white",
  legend_breaks = c(0, 1),
  col = circlize::colorRamp2(legend_breaks, c(min_colour, max_colour)),
  heatmap_legend_param = list(color_bar = "continuous", title = "p-value", at = c(0, 1)),
  rect_gp = grid::gpar(col = "black"),
  column_split_vector = NULL,
  row_split_vector = NULL,
  column_split = NULL,
  row_split = NULL,
  ...
)

```

**Arguments**

<code>ext_sol_df</code>	An <code>ext_solutions_df</code> class object (produced from the function <code>extend_solutions</code> ).
<code>order</code>	Numeric vector containing row order of the heatmap.
<code>cluster_columns</code>	Whether columns should be sorted by hierarchical clustering.
<code>cluster_rows</code>	Whether rows should be sorted by hierarchical clustering.
<code>show_row_names</code>	Whether row names should be shown.
<code>show_column_names</code>	Whether column names should be shown.
<code>min_colour</code>	Colour used for the lowest value in the heatmap.
<code>max_colour</code>	Colour used for the highest value in the heatmap.
<code>legend_breaks</code>	Numeric vector of breaks for the legend.
<code>col</code>	Colour function for <code>ComplexHeatmap::Heatmap()</code>
<code>heatmap_legend_param</code>	Legend function for <code>ComplexHeatmap::Heatmap()</code>
<code>rect_gp</code>	Cell border function for <code>ComplexHeatmap::Heatmap()</code>
<code>column_split_vector</code>	Vector of indices to split columns by.
<code>row_split_vector</code>	Vector of indices to split rows by.
<code>column_split</code>	Standard parameter of <code>ComplexHeatmap::Heatmap</code> .
<code>row_split</code>	Standard parameter of <code>ComplexHeatmap::Heatmap</code> .
<code>...</code>	Additional parameters passed to <code>ComplexHeatmap::Heatmap</code> .

**Value**

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the provided p-values.

**Examples**

```
#dl <- data_list(  
#   list(income, "household_income", "demographics", "ordinal"),  
#   list(pubertal, "pubertal_status", "demographics", "continuous"),  
#   list(fav_colour, "favourite_colour", "demographics", "categorical"),  
#   list(anxiety, "anxiety", "behaviour", "ordinal"),  
#   list(depress, "depressed", "behaviour", "ordinal"),  
#   uid = "unique_id"  
#)  
#  
#sc <- snf_config(  
#   dl,  
#   n_solutions = 4,  
#   dropout_dist = "uniform",  
#   max_k = 50  
#)  
#  
#sol_df <- batch_snf(dl, sc)  
#  
#ext_sol_df <- extend_solutions(sol_df, dl)  
#  
#pval_heatmap(ext_sol_df)
```

---

quality\_measures

*Quality metrics*

---

**Description**

These functions calculate conventional metrics of cluster solution quality.

**Usage**

```
calculate_silhouettes(sol_df)
```

```
calculate_dunn_indices(sol_df)
```

```
calculate_db_indices(sol_df)
```

**Arguments**

`sol_df` A `solutions_df` class object created by `batch_snf()` with the parameter `return_sim_mats = TRUE`.

## Details

`calculate_silhouettes`: A wrapper for `cluster::silhouette` that calculates silhouette scores for all cluster solutions in a provided solutions data frame. Silhouette values range from -1 to +1 and indicate an overall ratio of how close together observations within a cluster are to how far apart observations across clusters are. You can learn more about interpreting the results of this function by calling `?cluster::silhouette`.

`calculate_dunn_indices`: A wrapper for `clv::clv.Dunn` that calculates Dunn indices for all cluster solutions in a provided solutions data frame. Dunn indices, like silhouette scores, similarly reflect similarity within clusters and separation across clusters. You can learn more about interpreting the results of this function by calling `?clv::clv.Dunn`.

`calculate_db_indices`: A wrapper for `clv::clv.Davies.Bouldin` that calculates Davies-Bouldin indices for all cluster solutions in a provided solutions data frame. These values can be interpreted similarly as those above. You can learn more about interpreting the results of this function by calling `?clv::clv.Davies.Bouldin`.

## Value

A list of silhouette class objects, a vector of Dunn indices, or a vector of Davies-Bouldin indices depending on which function was used.

## Examples

```
input_dl <- data_list(
  list(gender_df, "gender", "demographics", "categorical"),
  list(diagnosis_df, "diagnosis", "clinical", "categorical"),
  uid = "patient_id"
)

sc <- snf_config(input_dl, n_solutions = 5)

sol_df <- batch_snf(input_dl, sc, return_sim_mats = TRUE)

# calculate Davies-Bouldin indices
davies_bouldin_indices <- calculate_db_indices(sol_df)

# calculate Dunn indices
dunn_indices <- calculate_dunn_indices(sol_df)

# calculate silhouette scores
silhouette_scores <- calculate_silhouettes(sol_df)
```

---

random\_removal

*Generate random removal sequence*

---

## Description

Helper function to contribute to rows within the settings data frame. Number of columns removed follows a uniform or exponential probability distribution.

**Usage**

```
random_removal(
  columns,
  min_removed_inputs,
  max_removed_inputs,
  dropout_dist = "exponential"
)
```

**Arguments**

`columns` Columns of the `settings_df` that are passed in

`min_removed_inputs` The smallest number of input data frames that may be randomly removed.

`max_removed_inputs` The largest number of input data frames that may be randomly removed.

`dropout_dist` Indication of how input data frames should be dropped. can be "none" (no dropout), "uniform" (uniformly draw number between min and max removed inputs), or "exponential" (like uniform, but using an exponential distribution; default).

**Value**

`inclusions_df` data frame that can be `rbind`'ed to the `settings_df`

---

`rbind.ext_solutions_df`

*Row-binding of solutions data frame class objects.*

---

**Description**

Row-binding of solutions data frame class objects.

**Usage**

```
## S3 method for class 'ext_solutions_df'
rbind(reset_indices = FALSE, ...)
```

**Arguments**

`reset_indices` If TRUE, re-labels the "solutions" indices in the solutions data frame from 1 to the number of defined settings.

`...` An arbitrary number of `ext_solutions_df` class objects.

**Value**

An `ext_solutions_df` class object.



---

rbind.solutions\_df      *Row-binding of solutions data frame class objects.*

---

**Description**

Row-binding of solutions data frame class objects.

**Usage**

```
## S3 method for class 'solutions_df'  
rbind(reset_indices = FALSE, ...)
```

**Arguments**

reset\_indices    If TRUE, re-labels the "solutions" indices in the solutions data frame from 1 to the number of defined settings.  
...              An arbitrary number of solutions\_df class objects.

**Value**

A solutions\_df class object.

---

rename\_dl              *Rename features in a data list*

---

**Description**

Rename features in a data list

**Usage**

```
rename_dl(dl, name_mapping)
```

**Arguments**

dl                  A nested list of input data from data\_list().  
name\_mapping      A named vector where the values are the features to be renamed and the names are the new names for those features.

**Value**

A data list ("list"-class object) with adjusted feature names.

## Examples

```
library(metasnf)

dl <- data_list(
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  list(anxiety, "anxiety", "behaviour", "ordinal"),
  list(depress, "depressed", "behaviour", "ordinal"),
  uid = "unique_id"
)

summary(dl, "feature")

name_changes <- c(
  "anxiety_score" = "cbcl_anxiety_r",
  "depression_score" = "cbcl_depress_r"
)

dl <- rename_dl(dl, name_changes)

summary(dl, "feature")
```

---

resample

*Helper resample function found in ?sample*

---

## Description

Like sample, but when given a single value x, returns back that single value instead of a random value from 1 to x.

## Usage

```
resample(x, ...)
```

## Arguments

x	Vector or single value to sample from
...	Remaining arguments for base::sample function

## Value

Numeric vector result of running base::sample.

---

save_heatmap	<i>Save a heatmap object to a file</i>
--------------	--

---

### Description

Save a heatmap object to a file

### Usage

```
save_heatmap(heatmap, path, width = 480, height = 480, res = 100)
```

### Arguments

heatmap	The heatmap object to save.
path	The path to save the heatmap to.
width	The width of the heatmap.
height	The height of the heatmap.
res	The resolution of the heatmap.

### Value

Does not return any value. Saves heatmap to file.

---

settings_df	<i>Build a settings data frame</i>
-------------	------------------------------------

---

### Description

The settings\_df is a data frame whose rows completely specify the hyperparameters and decisions required to transform individual input data frames (found in a data list, see ?data\_list) into a single similarity matrix through SNF. The format of the settings data frame is as follows:

- A column named "solution": This column is used to keep track of the rows and should have integer values only.
- A column named "alpha": This column contains the value of the alpha hyperparameter that will be used on that run of the SNF pipeline.
- A column named "k": Like above, but for the K (nearest neighbours) hyperparameter.
- A column named "t": Like above, but for the t (number of iterations) hyperparameter.
- A column named "snf\_scheme": Which of 3 pre-defined schemes will be used to integrate the data frames of the data list into a final fused network. The purpose of varying these schemes is primarily to increase the diversity of the generated cluster solutions.

- A value of 1 corresponds to the "individual" scheme, in which all data frames are directly merged by SNF into the final fused network. This scheme corresponds to the approach shown in the original SNF paper.
  - A value of 2 corresponds to the "two-step" scheme, in which all data frames within a domain are first merged into a domain-specific fused network. Next, domain-specific networks are fused once more by SNF into the final fused network. This scheme is useful for fairly re-weighting SNF pipelines with unequal numbers of data frames across domains.
  - A value of 3 corresponds to the "domain" scheme, in which all data frames within a domain are first concatenated into a single domain-specific data frame before being merged by SNF into the final fused network. This approach serves as an alternative way to re-weight SNF pipelines with unequal numbers of data frames across domains. You can learn more about this parameter here: [https://branchlab.github.io/metasnf/articles/snf\\_schemes.html](https://branchlab.github.io/metasnf/articles/snf_schemes.html).
- A column named "clust\_alg": Specification of which clustering algorithm will be applied to the final similarity matrix. By default, this column can take on the integer values 1 or 2, which correspond to spectral clustering where the number of clusters is determined by the eigen-gap or rotation cost heuristic respectively. You can learn more about this parameter here: [https://branchlab.github.io/metasnf/articles/clustering\\_algorithms.html](https://branchlab.github.io/metasnf/articles/clustering_algorithms.html).
  - A column named "cnt\_dist": Specification of which distance metric will be used for data frames of purely continuous data. You can learn about this metric and its defaults here: [https://branchlab.github.io/metasnf/articles/distance\\_metrics.html](https://branchlab.github.io/metasnf/articles/distance_metrics.html)
  - A column named "dsc\_dist": Like above, but for discrete data frames.
  - A column named "ord\_dist": Like above, but for ordinal data frames.
  - A column named "cat\_dist": Like above, but for categorical data frames.
  - A column named "mix\_dist": Like above, but for mixed-type (e.g., both categorical and discrete) data frames.
  - One column for every input data frame in the corresponding data list which can either have the value of 0 or 1. The name of the column should be formatted as "inc\_[]" where the square brackets are replaced with the name (as found in `dl_summary(dl)$"name"`) of each data frame. When 0, that data frame will be excluded from that run of the SNF pipeline. When 1, that data frame will be included.

## Usage

```
settings_df(
  dl,
  n_solutions = 0,
  min_removed_inputs = 0,
  max_removed_inputs = length(dl) - 1,
  dropout_dist = "exponential",
  min_alpha = NULL,
  max_alpha = NULL,
  min_k = NULL,
  max_k = NULL,
  min_t = NULL,
  max_t = NULL,
```

```

alpha_values = NULL,
k_values = NULL,
t_values = NULL,
possible_snf_schemes = c(1, 2, 3),
clustering_algorithms = NULL,
continuous_distances = NULL,
discrete_distances = NULL,
ordinal_distances = NULL,
categorical_distances = NULL,
mixed_distances = NULL,
dfl = NULL,
snf_input_weights = NULL,
snf_domain_weights = NULL,
retry_limit = 10,
allow_duplicates = FALSE
)

```

### Arguments

d1	A nested list of input data from <code>data_list()</code> .
n_solutions	Number of rows to generate for the settings data frame.
min_removed_inputs	The smallest number of input data frames that may be randomly removed. By default, 0.
max_removed_inputs	The largest number of input data frames that may be randomly removed. By default, this is 1 less than all the provided input data frames in the data list.
dropout_dist	Parameter controlling how the random removal of input data frames should occur. Can be "none" (no input data frames are randomly removed), "uniform" (uniformly sample between <code>min_removed_inputs</code> and <code>max_removed_inputs</code> to determine number of input data frames to remove), or "exponential" (pick number of input data frames to remove by sampling from <code>min_removed_inputs</code> to <code>max_removed_inputs</code> with an exponential distribution; the default).
min_alpha	The minimum value that the alpha hyperparameter can have. Random assigned value of alpha for each row will be obtained by uniformly sampling numbers between <code>min_alpha</code> and <code>max_alpha</code> at intervals of 0.1. Cannot be used in conjunction with the <code>alpha_values</code> parameter.
max_alpha	The maximum value that the alpha hyperparameter can have. See <code>min_alpha</code> parameter. Cannot be used in conjunction with the <code>alpha_values</code> parameter.
min_k	The minimum value that the k hyperparameter can have. Random assigned value of k for each row will be obtained by uniformly sampling numbers between <code>min_k</code> and <code>max_k</code> at intervals of 1. Cannot be used in conjunction with the <code>k_values</code> parameter.
max_k	The maximum value that the k hyperparameter can have. See <code>min_k</code> parameter. Cannot be used in conjunction with the <code>k_values</code> parameter.
min_t	The minimum value that the t hyperparameter can have. Random assigned value of t for each row will be obtained by uniformly sampling numbers between

	min_t and max_t at intervals of 1. Cannot be used in conjunction with the t_values parameter.
max_t	The maximum value that the t hyperparameter can have. See min_t parameter. Cannot be used in conjunction with the t_values parameter.
alpha_values	A number or numeric vector of a set of possible values that alpha can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_alpha or max_alpha parameters.
k_values	A number or numeric vector of a set of possible values that k can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_k or max_k parameters.
t_values	A number or numeric vector of a set of possible values that t can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_t or max_t parameters.
possible_snf_schemes	A vector containing the possible snf_schemes to uniformly randomly select from. By default, the vector contains all 3 possible schemes: c(1, 2, 3). 1 corresponds to the "individual" scheme, 2 corresponds to the "domain" scheme, and 3 corresponds to the "twostep" scheme.
clustering_algorithms	A list of clustering algorithms to uniformly randomly pick from when clustering. When not specified, randomly select between spectral clustering using the eigen-gap heuristic and spectral clustering using the rotation cost heuristic. See ?clust_fns_list for more details on running custom clustering algorithms.
continuous_distances	A vector of continuous distance metrics to use when a custom dist_fns_list is provided.
discrete_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
ordinal_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
categorical_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
mixed_distances	A vector of mixed distance metrics to use when a custom dist_fns_list is provided.
df1	List containing distance metrics to vary over. See ?generate_dist_fns_list.
snf_input_weights	Nested list containing weights for when SNF is used to merge individual input measures (see ?generate_snf_weights)
snf_domain_weights	Nested list containing weights for when SNF is used to merge domains (see ?generate_snf_weights)

`retry_limit` The maximum number of attempts to generate a novel row. This function does not return matrices with identical rows. As the range of requested possible settings tightens and the number of requested rows increases, the risk of randomly generating a row that already exists increases. If a new random row has matched an existing row `retry_limit` number of times, the function will terminate.

`allow_duplicates` If TRUE, enables creation of a settings data frame with duplicate non-feature weighting related hyperparameters. This function should only be used when paired with a custom weights matrix that has non-duplicate rows.

**Value**

A settings data frame

---

shiny_annotator	<i>Launch a shiny app to identify meta cluster boundaries</i>
-----------------	---

---

**Description**

This function calls the `htShiny()` function from the package `InteractiveComplexHeatmap` to assist users in identifying the indices of the boundaries between meta clusters in a meta cluster heatmap. By providing a heatmap of inter-solution similarities (obtained through `meta_cluster_heatmap()`), users can click on positions within the heatmap that appear to meaningfully separate major sets of similar cluster solutions by visual inspection. The corresponding indices of the clicked positions are printed to the console and also shown within the app. This function can only run from an interactive session of R.

**Usage**

```
shiny_annotator(ari_heatmap)
```

**Arguments**

`ari_heatmap` Heatmap of ARIs to divide into meta clusters.

**Value**

Does not return any value. Launches interactive shiny applet.

**Examples**

```
#dl <- data_list(
#   list(cort_sa, "cortical_surface_area", "neuroimaging", "continuous"),
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
#)
```

```
#
#set.seed(42)
#my_sc <- snf_config(
#  dl = dl,
#  n_solutions = 20,
#  min_k = 20,
#  max_k = 50
#)
#
#sol_df <- batch_snf(dl, my_sc)
#
#sol_aris <- calc_aris(sol_df)
#
#meta_cluster_order <- get_matrix_order(sol_aris)
#
#ari_hm <- meta_cluster_heatmap(sol_aris, order = meta_cluster_order)
#
## Click on meta cluster boundaries to obtain `split_vec` values
#shiny_annotator(ari_hm)
#
#split_vec <- c(6, 10, 16)
#
#ari_hm <- meta_cluster_heatmap(
#  sol_aris,
#  order = meta_cluster_order,
#  split_vector = split_vec
#)
```

---

similarity\_matrix\_heatmap

*Plot heatmap of similarity matrix*

---

## Description

Plot heatmap of similarity matrix

## Usage

```
similarity_matrix_heatmap(
  similarity_matrix,
  order = NULL,
  cluster_solution = NULL,
  scale_diag = "mean",
  log_graph = TRUE,
  cluster_rows = FALSE,
  cluster_columns = FALSE,
  show_row_names = FALSE,
  show_column_names = FALSE,
  data = NULL,
```



```

    left_bar = NULL,
    right_bar = NULL,
    top_bar = NULL,
    bottom_bar = NULL,
    left_hm = NULL,
    right_hm = NULL,
    top_hm = NULL,
    bottom_hm = NULL,
    annotation_colours = NULL,
    min_colour = NULL,
    max_colour = NULL,
    split_vector = NULL,
    row_split = NULL,
    column_split = NULL,
    ...
)

```

### Arguments

similarity_matrix	A similarity matrix
order	Vector of numbers to reorder the similarity matrix (and data if provided). Overwrites ordering specified by cluster_solution param.
cluster_solution	Row of a solutions data frame or column of a transposed solutions data frame.
scale_diag	Method of rescaling matrix diagonals. Can be "none" (don't change diagonals), "mean" (replace diagonals with average value of off-diagonals), or "zero" (replace diagonals with 0).
log_graph	If TRUE, log transforms the graph.
cluster_rows	Parameter for ComplexHeatmap::Heatmap.
cluster_columns	Parameter for ComplexHeatmap::Heatmap.
show_row_names	Parameter for ComplexHeatmap::Heatmap.
show_column_names	Parameter for ComplexHeatmap::Heatmap.
data	A data frame containing elements requested for annotation.
left_bar	Named list of strings, where the strings are features in df that should be used for a barplot annotation on the left of the plot and the names are the names that will be used to caption the plots and their legends.
right_bar	See left_bar.
top_bar	See left_bar.
bottom_bar	See left_bar.
left_hm	Like left_bar, but with a heatmap annotation instead of a barplot annotation.
right_hm	See left_hm.

top_hm	See left_hm.
bottom_hm	See left_hm.
annotation_colours	Named list of heatmap annotations and their colours.
min_colour	Colour used for the lowest value in the heatmap.
max_colour	Colour used for the highest value in the heatmap.
split_vector	A vector of partition indices.
row_split	Standard parameter of ComplexHeatmap: :Heatmap.
column_split	Standard parameter of ComplexHeatmap: :Heatmap.
...	Additional parameters passed into ComplexHeatmap::Heatmap.

### Value

Returns a heatmap (class "Heatmap" from package ComplexHeatmap) that displays the similarities between observations in the provided matrix.

### Examples

```
#my_dl <- data_list(
#   list(
#     data = expression_df,
#     name = "expression_data",
#     domain = "gene_expression",
#     type = "continuous"
#   ),
#   list(
#     data = methylation_df,
#     name = "methylation_data",
#     domain = "gene_methylation",
#     type = "continuous"
#   ),
#   uid = "patient_id"
#)
#
#sc <- snf_config(my_dl, n_solutions = 10)
#
#sol_df <- batch_snf(my_dl, sc, return_sim_mats = TRUE)
#
#sim_mats <- sim_mats_list(sol_df)
#
#similarity_matrix_heatmap(
#   sim_mats[[1]],
#   cluster_solution = sol_df[1, ]
#)
```

---

sim_mats_list	<i>Create or extract a sim_mats_list class object</i>
---------------	---

---

**Description**

Create or extract a sim\_mats\_list class object

**Usage**

```
sim_mats_list(x)
```

**Arguments**

x                    The object to create or extract a sim\_mats\_list from.

**Value**

A sim\_mats\_list class object.

---

siw_euclidean_distance	<i>Squared (including weights) Euclidean distance</i>
------------------------	---

---

**Description**

Squared (including weights) Euclidean distance

**Usage**

```
siw_euclidean_distance(df, weights_row)
```

**Arguments**

df                    data frame containing at least 1 data column.  
weights\_row          Single-row data frame where the column names contain the column names in df  
and the row contains the corresponding weights.

**Value**

distance\_matrix A distance matrix.

---

snf_config	<i>Define configuration for generating a set of SNF-based cluster solutions</i>
------------	---

---

### Description

snf\_config() constructs an SNF config object which inherits from classes snf\_config and list. This object is used to store all settings required to transform data stored in a data\_list class object into a space of cluster solutions by SNF. The SNF config object contains the following components: 1. A settings data frame (inherits from settings\_df and data.frame). Data frame that stores SNF-specific hyperparameters and information about feature selection and weighting, SNF schemes, clustering algorithms, and distance metrics. Each row of the settings data frame corresponds to a distinct cluster solution. 2. A clustering algorithms list (inherits from clust\_fns\_list and list), which stores all clustering algorithms that the settings data frame can point to. 3. A distance metrics list (inherits from dist\_metrics\_list and list), which stores all distance metrics that the settings data frame can point to. 4. A weights matrix (inherits from weights\_matrix, matrix, and array'), which stores the feature weights to use prior to distance calculations. Each column of the weights matrix corresponds to a different feature in the data list and each row corresponds to a different row in the settings data frame.

### Usage

```
snf_config(
  dl = NULL,
  sdf = NULL,
  dfl = NULL,
  cfl = NULL,
  wm = NULL,
  n_solutions = 0,
  min_removed_inputs = 0,
  max_removed_inputs = length(dl) - 1,
  dropout_dist = "exponential",
  min_alpha = NULL,
  max_alpha = NULL,
  min_k = NULL,
  max_k = NULL,
  min_t = NULL,
  max_t = NULL,
  alpha_values = NULL,
  k_values = NULL,
  t_values = NULL,
  possible_snf_schemes = c(1, 2, 3),
  clustering_algorithms = NULL,
  continuous_distances = NULL,
  discrete_distances = NULL,
  ordinal_distances = NULL,
  categorical_distances = NULL,
```

```

mixed_distances = NULL,
snf_input_weights = NULL,
snf_domain_weights = NULL,
retry_limit = 10,
cnt_dist_fns = NULL,
dsc_dist_fns = NULL,
ord_dist_fns = NULL,
cat_dist_fns = NULL,
mix_dist_fns = NULL,
automatic_standard_normalize = FALSE,
use_default_dist_fns = FALSE,
clust_fns = NULL,
use_default_clust_fns = FALSE,
weights_fill = "ones"
)

```

### Arguments

<code>dl</code>	A nested list of input data from <code>data_list()</code> .
<code>sdf</code>	A <code>settings_df</code> class object. Overrides settings data frame related parameters.
<code>df1</code>	A <code>dist_fns_list</code> class object. Overrides distance functions list related parameters.
<code>cfl</code>	A <code>clust_fns_list</code> class object. Overrides clustering functions list related parameters.
<code>wm</code>	A <code>weights_matrix</code> class object. Overrides weights matrix related parameters.
<code>n_solutions</code>	Number of rows to generate for the settings data frame.
<code>min_removed_inputs</code>	The smallest number of input data frames that may be randomly removed. By default, 0.
<code>max_removed_inputs</code>	The largest number of input data frames that may be randomly removed. By default, this is 1 less than all the provided input data frames in the data list.
<code>dropout_dist</code>	Parameter controlling how the random removal of input data frames should occur. Can be "none" (no input data frames are randomly removed), "uniform" (uniformly sample between <code>min_removed_inputs</code> and <code>max_removed_inputs</code> to determine number of input data frames to remove), or "exponential" (pick number of input data frames to remove by sampling from <code>min_removed_inputs</code> to <code>max_removed_inputs</code> with an exponential distribution; the default).
<code>min_alpha</code>	The minimum value that the alpha hyperparameter can have. Random assigned value of alpha for each row will be obtained by uniformly sampling numbers between <code>min_alpha</code> and <code>max_alpha</code> at intervals of 0.1. Cannot be used in conjunction with the <code>alpha_values</code> parameter.
<code>max_alpha</code>	The maximum value that the alpha hyperparameter can have. See <code>min_alpha</code> parameter. Cannot be used in conjunction with the <code>alpha_values</code> parameter.
<code>min_k</code>	The minimum value that the k hyperparameter can have. Random assigned value of k for each row will be obtained by uniformly sampling numbers between

	min_k and max_k at intervals of 1. Cannot be used in conjunction with the k_values parameter.
max_k	The maximum value that the k hyperparameter can have. See min_k parameter. Cannot be used in conjunction with the k_values parameter.
min_t	The minimum value that the t hyperparameter can have. Random assigned value of t for each row will be obtained by uniformly sampling numbers between min_t and max_t at intervals of 1. Cannot be used in conjunction with the t_values parameter.
max_t	The maximum value that the t hyperparameter can have. See min_t parameter. Cannot be used in conjunction with the t_values parameter.
alpha_values	A number or numeric vector of a set of possible values that alpha can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_alpha or max_alpha parameters.
k_values	A number or numeric vector of a set of possible values that k can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_k or max_k parameters.
t_values	A number or numeric vector of a set of possible values that t can take on. Value will be obtained by uniformly sampling the vector. Cannot be used in conjunction with the min_t or max_t parameters.
possible_snf_schemes	A vector containing the possible snf_schemes to uniformly randomly select from. By default, the vector contains all 3 possible schemes: c(1, 2, 3). 1 corresponds to the "individual" scheme, 2 corresponds to the "domain" scheme, and 3 corresponds to the "twostep" scheme.
clustering_algorithms	A list of clustering algorithms to uniformly randomly pick from when clustering. When not specified, randomly select between spectral clustering using the eigen-gap heuristic and spectral clustering using the rotation cost heuristic. See ?clust_fns_list for more details on running custom clustering algorithms.
continuous_distances	A vector of continuous distance metrics to use when a custom dist_fns_list is provided.
discrete_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
ordinal_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
categorical_distances	A vector of categorical distance metrics to use when a custom dist_fns_list is provided.
mixed_distances	A vector of mixed distance metrics to use when a custom dist_fns_list is provided.

snf_input_weights	Nested list containing weights for when SNF is used to merge individual input measures (see ?generate_snf_weights)
snf_domain_weights	Nested list containing weights for when SNF is used to merge domains (see ?generate_snf_weights)
retry_limit	The maximum number of attempts to generate a novel row. This function does not return matrices with identical rows. As the range of requested possible settings tightens and the number of requested rows increases, the risk of randomly generating a row that already exists increases. If a new random row has matched an existing row <code>retry_limit</code> number of times, the function will terminate.
cnt_dist_fns	A named list of continuous distance metric functions.
dsc_dist_fns	A named list of discrete distance metric functions.
ord_dist_fns	A named list of ordinal distance metric functions.
cat_dist_fns	A named list of categorical distance metric functions.
mix_dist_fns	A named list of mixed distance metric functions.
automatic_standard_normalize	If TRUE, will automatically use standard normalization prior to calculation of any numeric distances. This parameter overrides all other distance functions list-related parameters.
use_default_dist_fns	If TRUE, prepend the base distance metrics (euclidean distance for continuous, discrete, and ordinal data and gower distance for categorical and mixed data) to the resulting distance metrics list.
clust_fns	A list of named clustering functions
use_default_clust_fns	If TRUE, prepend the base clustering algorithms ( <code>spectral_eigen</code> and <code>spectral_rot</code> , which apply spectral clustering and use the eigen-gap and rotation cost heuristics respectively for determining the number of clusters in the graph) to <code>clust_fns</code> .
weights_fill	String indicating what to populate generate rows with. Can be "ones" (default; fill matrix with 1), "uniform" (fill matrix with uniformly distributed random values), or "exponential" (fill matrix with exponentially distributed random values).

**Value**

An `snf_config` class object.

**Examples**

```
# Simple random config for 5 cluster solutions
input_dl <- data_list(
  list(anxiety, "anxiety", "behaviour", "ordinal"),
  list(depress, "depressed", "behaviour", "ordinal"),
  uid = "unique_id"
)
my_sc <- snf_config(
```

```

    dl = input_dl,
    n_solutions = 5
  )

# specifying possible K range
my_sc <- snf_config(
  dl = input_dl,
  n_solutions = 5,
  min_k = 20,
  max_k = 40
)

# Random feature weights across from uniform distribution
my_sc <- snf_config(
  dl = input_dl,
  n_solutions = 5,
  min_k = 20,
  max_k = 40,
  weights_fill = "uniform"
)

# Specifying custom pre-built clustering and distance functions
# - Random alternation between 2-cluster and 5-cluster solutions
# - When continuous or discrete data frames are being processed,
#   randomly alternate between standardized/normalized Euclidean
#   distance vs. regular Euclidean distance
my_sc <- snf_config(
  dl = input_dl,
  n_solutions = 5,
  min_k = 20,
  max_k = 40,
  weights_fill = "uniform",
  clust_fns = list(
    "two_cluster_spectral" = spectral_two,
    "five_cluster_spectral" = spectral_five
  ),
  cnt_dist_fns = list(
    "euclidean" = euclidean_distance,
    "std_nrm_euc" = sn_euclidean_distance
  ),
  dsc_dist_fns = list(
    "euclidean" = euclidean_distance,
    "std_nrm_euc" = sn_euclidean_distance
  )
)

```

---

split\_parser

*Helper function to determine which row and columns to split on*


---

### Description

Helper function to determine which row and columns to split on



**Usage**

```
split_parser(
  row_split_vector = NULL,
  column_split_vector = NULL,
  row_split = NULL,
  column_split = NULL,
  n_rows,
  n_columns
)
```

**Arguments**

`row_split_vector` A vector of row indices to split on.

`column_split_vector` A vector of column indices to split on.

`row_split` Standard parameter of `ComplexHeatmap::Heatmap`.

`column_split` Standard parameter of `ComplexHeatmap::Heatmap`.

`n_rows` The number of rows in the data.

`n_columns` The number of columns in the data.

**Value**

"list"-class object containing `row_split` and `column_split` character vectors to pass into `ComplexHeatmap::Heatmap`.

---

subc_v	<i>Mock ABCD subcortical volumes data</i>
--------	---

---

**Description**

Like the mock data frame "abcd\_subc\_v", but with "unique\_id" as the "uid".

**Usage**

```
subc_v
```

**Format**

`subc_v`:  
 A data frame with 174 rows and 31 columns:  
**unique\_id** The unique identifier of the ABCD dataset  
 ... Subcortical volumes of various ROIs (mm<sup>3</sup>, I think)

**Source**

Though this data is no longer "real" ABCD data, the reference for using ABCD as a data source is below:

Data used in the preparation of this article were obtained from the Adolescent Brain Cognitive DevelopmentSM (ABCD) Study (<https://abcdstudy.org>), held in the NIMH Data Archive (NDA). This is a multisite, longitudinal study designed to recruit more than 10,000 children age 9-10 and follow them over 10 years into early adulthood. The ABCD Study® is supported by the National Institutes of Health and additional federal partners under award numbers U01DA041048, U01DA050989, U01DA051016, U01DA041022, U01DA051018, U01DA051037, U01DA050987, U01DA041174, U01DA041106, U01DA041117, U01DA041028, U01DA041134, U01DA050988, U01DA051039, U01DA041156, U01DA041025, U01DA041120, U01DA051038, U01DA041148, U01DA041093, U01DA041089, U24DA041123, U24DA041147. A full list of supporters is available at <https://abcdstudy.org/federal-partners.html>. A listing of participating sites and a complete listing of the study investigators can be found at [https://abcdstudy.org/consortium\\_members/](https://abcdstudy.org/consortium_members/). ABCD consortium investigators designed and implemented the study and/or provided data but did not necessarily participate in the analysis or writing of this report. This manuscript reflects the views of the authors and may not reflect the opinions or views of the NIH or ABCD consortium investigators.

---

subsample\_dl

*Create subsamples of a data list*


---

**Description**

Given a data list, return a list of smaller data lists that are generated through random sampling (without replacement). The results of this function can be passed into `batch_snf_subsamples()` to obtain a list of resampled solutions data frames.

**Usage**

```
subsample_dl(
  dl,
  n_subsamples,
  subsample_fraction = NULL,
  n_observations = NULL
)
```

**Arguments**

`dl` A nested list of input data from `data_list()`.

`n_subsamples` Number of subsamples to create.

`subsample_fraction` Percentage of patients to include per subsample.

`n_observations` Number of patients to include per subsample.

**Value**

A "list" class object containing `n_subsamples` number of data lists. Each of those data lists contains a random `subsample_fraction` fraction of the observations of the provided data list.

**Examples**

```
# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
# )
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
```

---

subsample\_pairwise\_aris

*Calculate pairwise adjusted Rand indices across subsamples of data*

---

**Description**

Given a list of subsampled solutions data frames from `batch_snf_subsamples()`, this function calculates the adjusted Rand indices across all the subsamples of each solution. ARI calculation between two subsamples only factors in observations that were present in both subsamples.

**Usage**

```
subsample_pairwise_aris(subsample_solutions, verbose = FALSE)
```

**Arguments**

`subsample_solutions`

A list of solutions data frames from subsamples of the data. This object is generated by the function `batch_snf_subsamples()`.

`verbose`

If TRUE, output progress to console.

**Value**

A two-item list: "raw\_aris", a list of inter-subsample pairwise ARI matrices (one for each full cluster solution) and "ari\_summary", a data frame containing the mean and SD of the inter-subsample ARIs for each original cluster solution.

**Examples**

```

# my_dl <- data_list(
#   list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
#   list(income, "household_income", "demographics", "continuous"),
#   list(pubertal, "pubertal_status", "demographics", "continuous"),
#   uid = "unique_id"
# )
#
# sc <- snf_config(my_dl, n_solutions = 5, max_k = 40)
#
# my_dl_subsamples <- subsample_dl(
#   my_dl,
#   n_subsamples = 20,
#   subsample_fraction = 0.85
# )
#
# batch_subsample_results <- batch_snf_subsamples(
#   my_dl_subsamples,
#   sc,
#   verbose = TRUE
# )
#
# pairwise_aris <- subsample_pairwise_aris(
#   batch_subsample_results
#   verbose = TRUE
# )
#
# # Visualize ARIs
# ComplexHeatmap::Heatmap(
#   pairwise_aris[[1]],
#   heatmap_legend_param = list(
#     color_bar = "continuous",
#     title = "Inter-Subsample\nARI",
#     at = c(0, 0.5, 1)
#   ),
#   show_column_names = FALSE,
#   show_row_names = FALSE
# )

```

---

summarize\_clust\_fns\_list

*Summarize a clust\_fns\_list object*

---

**Description**

Summarize a clust\_fns\_list object

**Usage**

```
summarize_clust_fns_list(cfl)
```

**Arguments**

`cfl` A `clust_fns_list` class object.

**Value**

`summary_df` "data.frame" class object containing the name and index of each clustering algorithm in te provided `clust_fns_list`.

---

<code>summarize_dfl</code>	<i>Summarize metrics contained in a <code>dist_fns_list</code></i>
----------------------------	--

---

**Description**

Summarize metrics contained in a `dist_fns_list`

**Usage**

```
summarize_dfl(dist_fns_list)
```

**Arguments**

`dist_fns_list` A `dist_fns_list`.

**Value**

"data.frame"-class object summarizing items in a distance metrics list.

---

<code>summarize_dl</code>	<i>Summarize a data list</i>
---------------------------	------------------------------

---

**Description**

**[Deprecated]** Defunct function for summarizing a data list. Please use `summary()` instead.

**Usage**

```
summarize_dl(data_list, scope = "component")
```

**Arguments**

`data_list` A nested list of input data from `data_list()`.

`scope` The level of detail for the summary. Options are:

- "component" (default): One row per component (data frame) in the data list.
- "feature": One row for each feature in the data list.

**Value**

data.frame class object summarizing all components (or features if scope == "component").

---

summary.data\_list      *Summary method for class data\_list*

---

**Description**

Returns a data list summary (data.frame class object) containing information on components, features, variable types, domains, and component dimensions.

**Usage**

```
## S3 method for class 'data_list'
summary(object, scope = "component", ...)
```

**Arguments**

object	A data_list class object.
scope	The level of detail for the summary. By default, this is set to "component", which returns a summary of the data list at the component level. Can also be set to "feature", resulting in a summary at the feature level.
...	Other arguments passed to summary (not used in this function)

**Value**

A data.frame class object. If scope is "component", each row shows the name, variable type, domain, and dimensions of each component. If scope is "feature", each row shows the name, variable type, and domain of each feature.

---

summary\_features      *Pull features used to calculate summary p-values from an object*

---

**Description**

Pull features used to calculate summary p-values from an object

**Usage**

```
summary_features(x)
```

**Arguments**

x	The object to extract summary features from.
---	--

**Value**

A character vector of summary features.

---

train_test_assign	<i>Training and testing split</i>
-------------------	-----------------------------------

---

**Description**

Given a vector of `uid_id` and a threshold, returns a list of which members should be in the training set and which should be in the testing set. The function relies on whether or not the absolute value of the Jenkins's `one_at_a_time` hash function exceeds the maximum possible value (2147483647) multiplied by the threshold.

**Usage**

```
train_test_assign(train_frac, uids, seed = 42)
```

**Arguments**

<code>train_frac</code>	The fraction (0 to 1) of observations for training
<code>uids</code>	A character vector of UIDs to be distributed into training and test sets.
<code>seed</code>	Seed used for Jenkins's <code>one_at_a_time</code> hash function.

**Value**

A named list containing the training and testing `uid_ids`.

---

<code>uids</code>	<i>Pull UIDs from an object</i>
-------------------	---------------------------------

---

**Description**

Pull UIDs from an object

**Usage**

```
uids(x)
```

**Arguments**

<code>x</code>	The object to extract UIDs from.
----------------	----------------------------------

**Value**

A character vector of UIDs.

---

validate\_solutions\_df *Validator for solutions\_df class object*

---

### Description

Validator for solutions\_df class object

### Usage

```
validate_solutions_df(sol_df1)
```

### Arguments

sol\_df1            A solutions data frame-like object to be validated and converted into a solutions data frame.

### Value

If sol\_df1 has a valid structure for a solutions\_df class object, returns the input unchanged. Otherwise, raises an error.

---

var\_manhattan\_plot    *Manhattan plot of feature-feature association p-values*

---

### Description

Manhattan plot of feature-feature association p-values

### Usage

```
var_manhattan_plot(  
  dl,  
  key_var,  
  neg_log_pval_thresh = 5,  
  threshold = NULL,  
  point_size = 5,  
  text_size = 20,  
  plot_title = NULL,  
  hide_x_labels = FALSE,  
  bonferroni_line = FALSE  
)
```



**Arguments**

dl	List of data frames containing data information.
key_var	Feature for which the association p-values of all other features are plotted.
neg_log_pval_thresh	Threshold for negative log p-values.
threshold	p-value threshold to plot dashed line at.
point_size	Size of points in the plot.
text_size	Size of text in the plot.
plot_title	Title of the plot.
hide_x_labels	If TRUE, hides x-axis labels.
bonferroni_line	If TRUE, plots a dashed black line at the Bonferroni-corrected equivalent of the p-value threshold.

**Value**

A Manhattan plot (class "gg", "ggplot") showing the association p-values of features against one key feature in a data list.

**Examples**

```
dl <- data_list(
  list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
  list(income, "household_income", "demographics", "continuous"),
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  list(anxiety, "anxiety", "behaviour", "ordinal"),
  list(depress, "depressed", "behaviour", "ordinal"),
  uid = "unique_id"
)

var_manhattan <- var_manhattan_plot(
  dl,
  key_var = "household_income",
  plot_title = "Correlation of Features with Household Income",
  text_size = 16,
  neg_log_pval_thresh = 3,
  threshold = 0.05
)
```

**Description**

Function for building a weights matrix independently of an SNF config. The weights matrix contains one row corresponding to each row of the settings data frame in an SNF config (one row for each resulting cluster solution) and one column for each feature in the data list used for clustering. Values of the weights matrix are passed to distance metrics functions during the conversion of input data frames to distance matrices. Typically, there is no need to use this function directly. Instead, users should provide weights matrix-building parameters to the `snf_config()` function.

**Usage**

```
weights_matrix(dl = NULL, n_solutions = 1, weights_fill = "ones")
```

**Arguments**

<code>dl</code>	A nested list of input data from <code>data_list()</code> .
<code>n_solutions</code>	Number of rows to generate the template weights matrix for.
<code>weights_fill</code>	String indicating what to populate generate rows with. Can be "ones" (default; fill matrix with 1), "uniform" (fill matrix with uniformly distributed random values), or "exponential" (fill matrix with exponentially distributed random values).

**Value**

`wm` A properly formatted matrix containing columns for all the features that require weights and rows.

**Examples**

```
input_dl <- data_list(
  list(subc_v, "subcortical_volume", "neuroimaging", "continuous"),
  list(income, "income", "demographics", "continuous"),
  list(pubertal, "pubertal_status", "demographics", "continuous"),
  uid = "unique_id"
)

sc <- snf_config(input_dl, n_solutions = 5)

wm <- weights_matrix(input_dl, n_solutions = 5, weights_fill = "uniform")

# updating an SNF config in parts
sc$weights_matrix <- wm
```

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