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# **data\_tools Documentation**

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# CONTENTS

<b>1</b>	<b>Disclaimer</b>	<b>3</b>
<b>2</b>	<b>Dependencies</b>	<b>5</b>
<b>3</b>	<b>Installation</b>	<b>7</b>
<b>4</b>	<b>Modules</b>	<b>9</b>
4.1	data_tools.databases . . . . .	9
4.2	data_tools.diffusion . . . . .	11
4.3	data_tools.models . . . . .	11
4.4	data_tools.plots . . . . .	13
4.5	data_tools.sets . . . . .	17
4.6	data_tools.strings . . . . .	19
	<b>Python Module Index</b>	<b>21</b>
	<b>Index</b>	<b>23</b>



Data tools: a collection of Python functions and classes designed to make data scientists' life easier.

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## DISCLAIMER

This package is still under development and will be periodically updated with new features. Contributions are very welcome (fork + pull request). If you find any bug or suggestion for upgrades, please use the [issue system](#).

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## DEPENDENCIES

- NumPy
- Matplotlib
- Pandas
- SciPy
- Scikit-learn



## INSTALLATION

First download/clone `data_tools` from the [GitHub repository](https://github.com/Nic-Nic/data_tools). From the terminal:

```
git clone https://github.com/Nic-Nic/data_tools.git
cd data_tools
```

Then you can install it by running `setup.py` as follows:

```
python setup.py sdist
```

Or using `pip`:

```
pip install .
```

Along with `data_tools`, all dependencies will be installed as well as the testing suite. In order to run the tests, type on the terminal:

```
python -m test_data_tools
```

**NOTE:** `data_tools.plots` module does not have any tests implemented.



## MODULES

### 4.1 data\_tools.databases

Databases functions module.

`data_tools.databases.kegg_link(query, target='pathway')`

Queries a request to the KEGG database to find related entries using cross-references. A list of available database(s) and query examples can be found in <https://www.kegg.jp/kegg/rest/keggapi.html#link>.

- **Arguments:**

- *query* [list]: Or any iterable type containing the identifier(s) to be queried as [str]. These can be either valid database identifiers or databases *per se* (see the link above).
- *target* [str]: Optional, 'pathway' by default. Targeted database to which the query should be linked to. You can check other options available in the URL above.

- **Returns:**

- [pandas.DataFrame]: Two-column table containing both the input query identifiers and their linked ones.

- **Example:**

```
>>> my_query = ['hsa:10458', 'ece:Z5100']
>>> kegg_link(my_query)
  query      pathway
0  hsa:10458  path:hsa04520
1  hsa:10458  path:hsa04810
2  ece:Z5100  path:ece05130
```

`data_tools.databases.kegg_pathway_mapping(df, mapid, filename=None)`

Makes a request to KEGG pathway mapping tool according to a given pathway ID (see [https://www.kegg.jp/kegg/tool/map\\_pathway2.html](https://www.kegg.jp/kegg/tool/map_pathway2.html) for more information). The user must provide a query of IDs to be mapped with their corresponding background colors (and optionally also foreground colors). The result is downloaded in the current directory or a user-specified path.

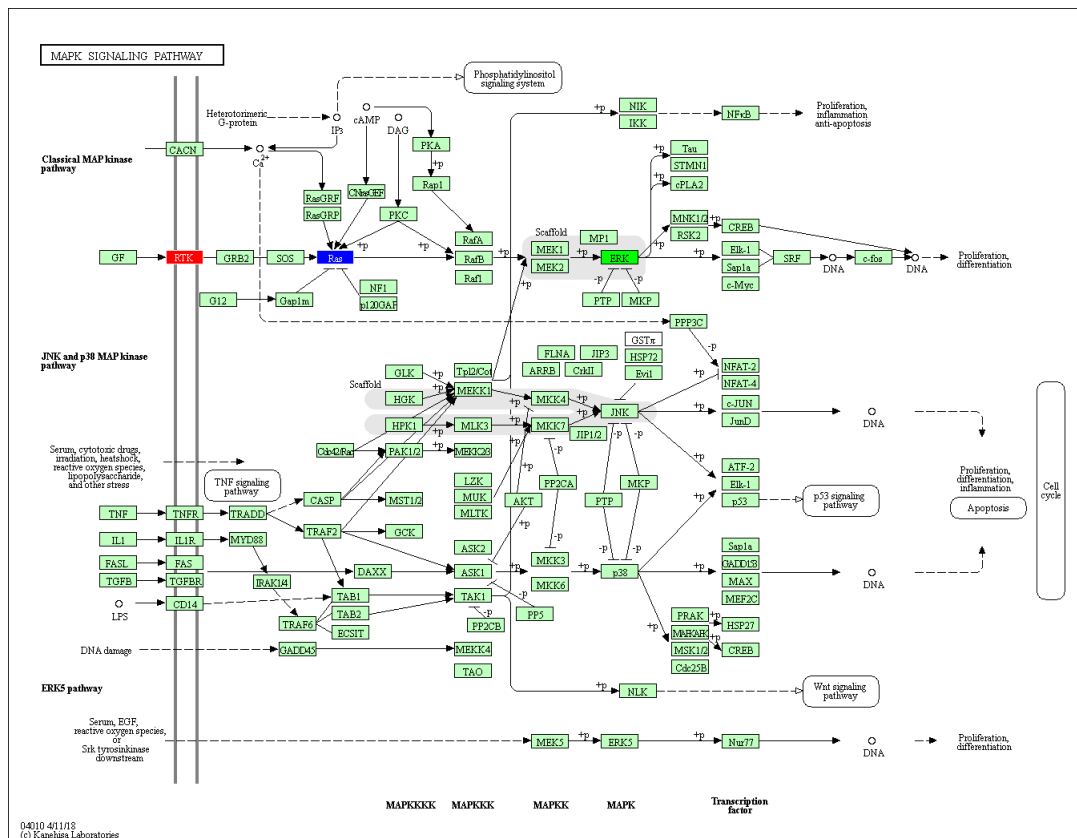
- **Arguments:**

- *df* [pandas.DataFrame]: Dataframe containing KEGG valid IDs in the first column and corresponding background colors (e.g.: red, blue, ...). Optionally, a third column with the foreground (font) colors can also be provided (black by default). **NOTE:** hexadecimal codes for colors is also supported. Index and column names of dataframe are ignored.
- *mapid* [str]: A valid KEGG pathway ID. It can be a general (e.g.: “mapXXXXX”) or organism-specific ID (e.g.: “hsaXXXXX”).

- *filename* [str]: Optional, *None* by default. This is, the image will be stored in the current directory with the *mapid* provided as file name. If provided, the image will be stored within the specified path/file name.

- **Example:**

```
>>> my_query = pandas.DataFrame([[ '1956', 'red', '#f1f1f1'],
...                               [ '3845', 'blue', '#f1f1f1'],
...                               [ '5594', 'green', 'black']])
>>> kegg_pathway_mapping(my_query, 'hsa04010')
```



`data_tools.databases.up_map(query, source='ACC', target='GENENAME')`

Queries a request to UniProt.org in order to map a given list of identifiers. You can check the options available of input/output identifiers at [https://www.uniprot.org/help/api\\_idmapping](https://www.uniprot.org/help/api_idmapping).

- **Arguments:**

- *query* [list]: Or any iterable type containing the identifiers to be queried as [str].
- *source* [str]: Optional, 'ACC' by default. This is, UniProt accession number. You can check other options available in the URL above.
- *target* [str]: Optional, 'GENENAME' by default. You can check other options available in the URL above.

- **Returns:**

- [pandas.DataFrame]: Two-column table containing both the inputted identifiers and the mapping result of these.

- **Examples:**

```

>>> my_query = ['P00533', 'P31749', 'P16220']
>>> up_map(my_query)
      ACC GENENAME
0  P00533    EGFR
1  P31749    AKT1
2  P16220    CREB1
>>> up_map(my_query, target='KEGG_ID')
      ACC    KEGG_ID
0  P00533  hsa:1956
2  P16220  hsa:1385
1  P31749  hsa:207

```

## 4.2 data\_tools.diffusion

Diffusion solvers module.

`data_tools.diffusion.euler_explicit1D(x0, dt, dx2, d=1, bcs='periodic')`

Computes diffusion on a 1D space over a time-step using Euler explicit method.

- **Arguments:**

- `x0` [numpy.ndarray]: Initial state of a 1D array from which the diffusion is to be computed.
- `dt` [float]: Discretization time-step.
- `dx2` [float]: Discretization spatial-step (squared).
- `d` [float]: Diffusion coefficient.
- `bcs` [str]: Optional, 'periodic' by default. Determines the boundary conditions. Available options are 'periodic', 'dirichlet' or 'neumann'.

- **Returns:**

- [numpy.ndarray]: Computed state array (1D) after one time-step according to the parameters and conditions selected.

## 4.3 data\_tools.models

Model classes module.

**class** `data_tools.models.DoseResponse(d_data, r_data, x0=[1, 1, 1], x_scale=[1, 1, 1], bounds=([0, 0, -inf], [inf, inf, inf]))`

Wrapper class for `scipy.optimize.least_squares` to fit dose-response curves on a pre-defined Hill function with the following form:

$$R = \frac{mD^n}{k^n + D^n}$$

Where  $D$  is the dose,  $k$ ,  $m$  and  $n$  are the parameters to be fitted.

- **Arguments:**

- `d_data` [numpy.ndarray]: Or any iterable (1D). Contains the training data corresponding to the dose.
- `r_data` [numpy.ndarray]: Or any iterable (1D). Contains the training data corresponding to the response.

- *x0* [list]: Optional, [1, 1, 1] by default. Or any iterable of three elements. Contains the initial guess for the parameters. Parameters are considered to be in alphabetical order. This is, first element corresponds to  $k$ , second is  $m$  and last is  $n$ .
- *x\_scale* [list]: Optional, [1, 1, 1] by default. Or any iterable of three elements. Scale of each parameter. May improve the fitting if the scaled parameters have similar effect on the cost function.
- *bounds* [tuple]: Optional ([0, 0, -np.inf], [np.inf, np.inf, np.inf]) by default. Two-element tuple containing the lower and upper boundaries for the parameters (elements of the tuple are iterables of three elements each).

- **Attributes:**

- *model* [scipy.optimize.OptimizeResult]: Contains the result of the optimized model. See [SciPy's reference](#) for more information.
- *params* [numpy.ndarray]: Three-element array containing the fitted parameters  $k$ ,  $m$  and  $n$ .

**ec** (*p*=50)

Computes the effective concentration for the specified percentage of maximal concentration ( $EC_p$ ).

- **Arguments:**

- *p* [int]: Optional, 50 by default ( $EC_{50}$ ). Defines the percentage of the maximal from which the effective concentration is to be computed.

- **Returns**

- [float]: Value of the  $EC_p$  computed according to the model parameters.

**plot** (*title*=None, *filename*=None, *figsize*=None, *legend*=True)

Plots the data points and the fitted function together.

- **Arguments:**

- *title* [str]: Optional, None by default. Defines the plot title.
- *filename* [str]: Optional, None by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- *figsize* [tuple]: Optional, None by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].
- *legend* [bool]: Optional, True by default. Indicates whether to show the plot legend or not.

- **Returns:**

- [matplotlib.figure.Figure]: Figure object showing the data points and the fitted model function.

**class** data\_tools.models.Lasso (*Cs*=500, *cv*=10, *sampler*='skf', *solver*='liblinear', \*\*kwargs)

Wrapper class inheriting from `sklearn.linear_model.LogisticRegressionCV` with L1 regularization.

- **Arguments:**

- *Cs* [int]: Optional, 500 by default. Integer or list of float values of regularization parameters to test. If an integer is passed, it will determine the number of values taken from a logarithmic scale between  $1e-4$  and  $1e4$ . Note that the value of the parameter is defined as the inverse of the regularization strength.
- *cv* [int]: Optional, 10 by default. Denotes the number of cross validation (CV) folds.



- *sampler* [str]: Optional, 'skf' by default. Determines which sampling method is used to generate the test and training sets for CV. Methods available are K-Fold ('kf'), Shuffle Split ('ss') and their stratified variants ('skf' and 'sss' respectively).
- *solver* [str]: Optional, 'liblinear' by default. Determines which solver algorithm to use. Note that L1 regularization can only be handled by 'liblinear' and 'saga'. Additionally if the classification is multinomial, only the latter option is available.
- *\*\*kwargs*: Optional. Any other keyword argument accepted by the `sklearn.linear_model.LogisticRegressionCV` class.

Other keyword arguments and functions available from the parent class `LogisticRegressionCV` can be found in [Scikit-Learn's reference](#).

**fit\_data** (*x*, *y*, *silent=False*)

Fits the data to the logistic model.

• **Arguments:**

- *x* [pandas.DataFrame]: Contains the values/measurements [float] of the features (columns) for each sample/replicate (rows).
- *y* [pandas.Series]: List or any iterable containing the observed class of each sample (must have the same order as in *x*).
- *silent* [bool]: Optional, `False` by default. Determines whether messages are printed or not.

**plot\_coef** (*filename=None*, *figsize=None*)

Plots the non-zero coefficients for the fitted predictor features.

• **Arguments:**

- *filename* [str]: Optional, `None` by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- *figsize* [tuple]: Optional, `None` by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

• **Returns:**

- [matplotlib.figure.Figure]: Figure object containing the bar plot of the non-zero coefficients.

**plot\_score** (*filename=None*, *figsize=None*)

Plots the mean score across all folds obtained during CV. The optimum C parameter chosen and its score are highlighted.

• **Arguments:**

- *filename* [str]: Optional, `None` by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- *figsize* [tuple]: Optional, `None` by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

• **Returns:**

- [matplotlib.figure.Figure]: Figure object containing the score plot.

## 4.4 data\_tools.plots

Plotting functions module.

`data_tools.plots.density(df, cvf=0.25, title=None, filename=None, figsize=None)`

Generates a density plot of the values on a data frame (row-wise).

- **Arguments:**

- `df` [pandas.DataFrame]: Contains the values to generate the plot. Each row is considered as an individual sample while each column contains a measured value.
- `cvf` [float]: Optional, 0.25 by default. Co-variance factor used in the gaussian kernel estimation. A higher value increases the smoothness.
- `title` [str]: Optional, `None` by default. Defines the plot title.
- `filename` [str]: Optional, `None` by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- `figsize` [tuple]: Optional, `None` by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

- **Returns:**

- [`matplotlib.figure.Figure`]: the figure object containing the density plot.

`data_tools.plots.piano_consensus(df, nchar=40, boxes=True, title=None, filename=None, figsize=None)`

Generates a GSEA consensus score plot like R package `piano`'s `consensusScores` function, but prettier. The main input is assumed to be a `pandas.DataFrame` whose data is the same as the `rankMat` from the result of `consensusScores`.

- **Arguments:**

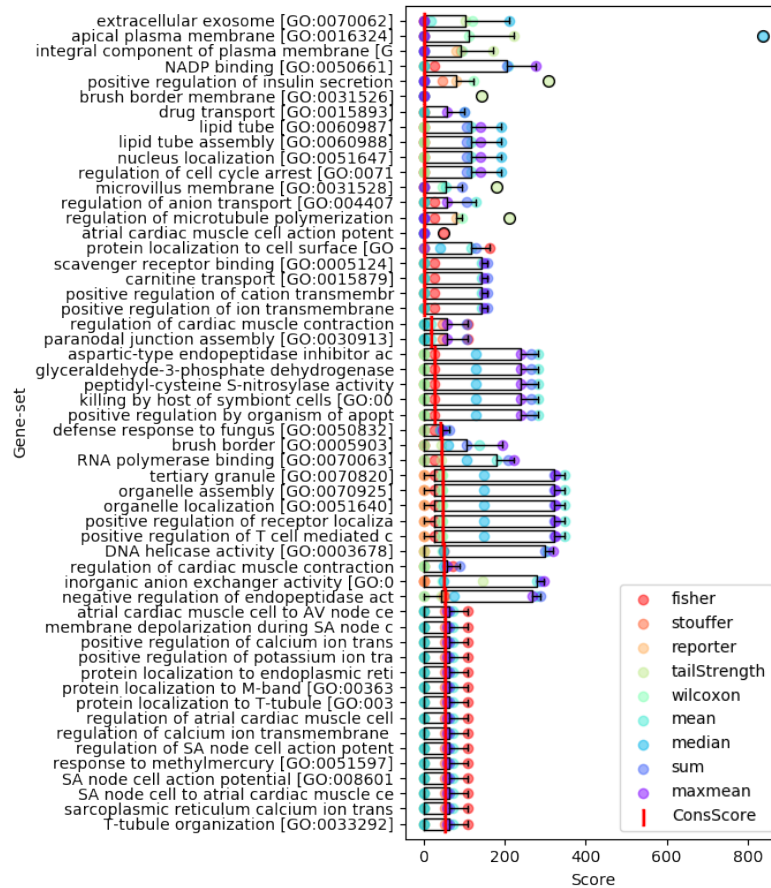
- `df` [pandas.DataFrame]: Values contained correspond to the scores of the gene-sets (consensus and each individual methods). Index must contain the gene-set labels. Columns are assumed to be `ConsRank` (ignored), `ConsScore` followed by the individual methods (e.g.: `mean`, `median`, `sum`, etc).
- `nchar` [int]: Optional, 40 by default. Number of string characters of the gene-set labels of the plot.
- `boxes` [bool]: Optional, `True` by default. Determines whether to show the boxplots of the gene-sets or not.
- `title` [str]: Optional, `None` by default. Defines the plot title.
- `filename` [str]: Optional, `None` by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- `figsize` [tuple]: Optional, `None` by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

- **Returns:**

- [`matplotlib.figure.Figure`]: the figure object containing a combination of box and scatter plots of the gene-set scores.

- **Example:**

```
>>> piano_consensus(df, figsize=[7, 8])
```



`data_tools.plots.venn(N, labels=['A', 'B', 'C', 'D', 'E'], c=['C0', 'C1', 'C2', 'C3', 'C4'], title=None, filename=None, figsize=None)`

Plots a Venn diagram from a list of sets  $N$ . Number of sets must be between 2 and 5 (inclusive).

- **Arguments:**

- $N$  [list]: Or any iterable type containing [set] objects.
- *labels* [list]: Optional, ['A', 'B', 'C', 'D', 'E'] by default. Labels for the sets following the same order as provided in  $N$ .
- *c* [list]: Optional, ['C0', 'C1', 'C2', 'C3', 'C4'] by default (matplotlib default colors). Any iterable containing color arguments tolerated by matplotlib (e.g.: ['r', 'b'] for red and blue). Must contain at least the same number of elements as  $N$  (if more are provided, they will be ignored).
- *title* [str]: Optional, None by default. Defines the plot title.
- *filename* [str]: Optional, None by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- *figsize* [tuple]: Optional, None by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

- **Returns:**

- [*matplotlib.figure.Figure*]: the figure object containing a combination of box and scatter plots of the gene-set scores.

- **Example:**

```
>>> N = [{0, 1}, {2, 3}, {1, 3, 4}] # Sets A, B, C
>>> venn(N)
```



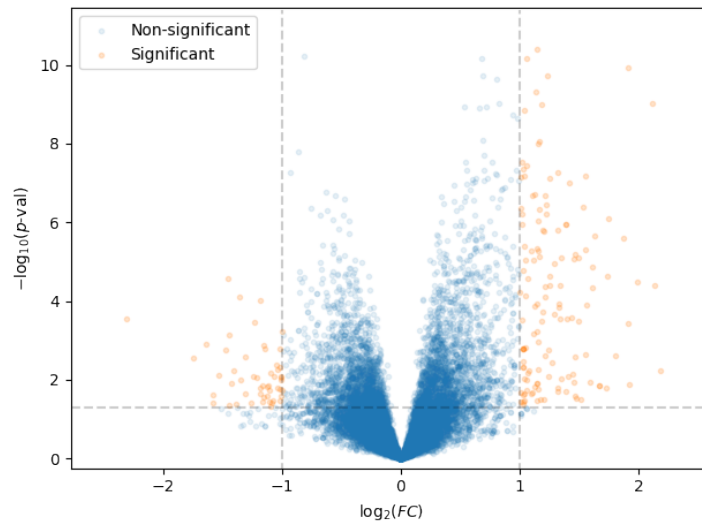
`data_tools.plots.volcano` (*logfc*, *logpval*, *thr\_pval*=0.05, *thr\_fc*=2.0, *c*=('C0', 'C1'), *legend*=True, *title*=None, *filename*=None, *figsize*=None)  
Generates a volcano plot from the differential expression data provided.

- **Arguments:**

- *logfc* [list]: Or any iterable type. Contains the log (usually base 2) fold-change values. Must have the same length as *logpval*.
- *logpval* [list]: Or any iterable type. Contains the -log p-values (usually base 10). Must have the same length as *logfc*.
- *thr\_pval* [float]: Optional, 0.05 by default. Specifies the p-value (non log-transformed) threshold to consider a measurement as significantly differentially expressed.
- *thr\_fc* [float]: Optional, 2. by default. Specifies the FC (non log-transformed) threshold to consider a measurement as significantly differentially expressed.
- *c* [tuple]: Optional, ('C0', 'C1') by default (matplotlib default colors). Any iterable containing two color arguments tolerated by matplotlib (e.g.: ['r', 'b'] for red and blue). First one is used for non-significant points, second for the significant ones.
- *legend* [bool]: Optional, True by default. Indicates whether to show the plot legend or not.
- *title* [str]: Optional, None by default. Defines the plot title.
- *filename* [str]: Optional, None by default. If passed, indicates the file name or path where to store the figure. Format must be specified (e.g.: .png, .pdf, etc)
- *figsize* [tuple]: Optional, None by default (default matplotlib size). Any iterable containing two values denoting the figure size (in inches) as [width, height].

- **Returns:**
  - [matplotlib.figure.Figure]: Figure object containing the volcano plot.
- **Example:**

```
>>> volcano(my_log_fc, my_log_pval)
```



## 4.5 data\_tools.sets

Set operations module.

`data_tools.sets.bit_or(a, b)`

Returns the bit operation OR between two bit-strings *a* and *b*. **NOTE:** *a* and *b* must have the same size.

- **Arguments:**
  - *a* [tuple]: Or any iterable type.
  - *b* [tuple]: Or any iterable type.
- **Returns:**
  - [tuple]: OR operation between *a* and *b* element-wise.
- **Examples:**

```
>>> a, b = (0, 0, 1), (1, 0, 1)
>>> bit_or(a, b)
(1, 0, 1)
```

`data_tools.sets.chunk_this(L, n)`

For a given list *L*, returns another list of *n*-sized chunks from it (in the same order).

- **Arguments:**
  - *L* [list]: The list to be sliced into sublists of the defined size.
  - *n* [int]: The size of the chunks.
- **Returns:**

- [list]: List of  $n$ -sized chunks from  $L$ . **NOTE:** If the number of items in  $L$  is not divisible by  $n$ , the last element returned will have an inferior size.

- **Examples:**

```
>>> L = range(6)
>>> chunk_this(L, 2)
[[0, 1], [2, 3], [4, 5]]
>>> chunk_this(L, 4)
[[0, 1, 2, 3], [4, 5]]
```

`data_tools.sets.find_min(A)`

Finds and returns the subset of vectors whose sum is minimum from a given set  $A$ .

- **Arguments:**

- $A$  [set]: Set of vectors ([tuple] or any iterable).

- **Returns:**

- [set]: Subset of vectors in  $A$  whose sum is minimum.

- **Examples:**

```
>>> A = {(0, 1, 1), (0, 1, 0), (1, 0, 0), (1, 1, 1)}
>>> find_min(A)
set([(0, 1, 0), (1, 0, 0)])
```

`data_tools.sets.in_all(x, N)`

Checks if a vector  $x$  is present in all sets contained in a list  $N$ .

- **Arguments:**

- $x$  [tuple]: Or any hashable type as long as is the same contained in the sets of  $N$ .
- $N$  [list]: Or any iterable type containing [set] objects.

- **Returns:**

- [bool]: True if  $x$  is found in all sets of  $N$ , False otherwise.

- **Examples:**

```
>>> N = [{(0, 0), (0, 1)}, # <- set A
...      {(0, 0), (1, 1), (1, 0)}] # <- set B
>>> x = (0, 0)
>>> in_all(x, N)
True
>>> y = (0, 1)
>>> in_all(y, N)
False
```

`data_tools.sets.subsets(N)`

Function that computes all possible logical relations between all sets on a list  $N$  and returns all subsets. This is, the subsets that would represent each intersecting area on a Venn diagram.

- **Arguments:**

- $N$  [list]: Or any iterable type containing [set] objects.

- **Returns:**

- [dict]: Collection of subsets according to the logical relations between the sets in  $N$ . The keys are binary codes that denote the logical relation (see examples below).

- **Examples:**

```
>>> N = [{0, 1, 2}, {2, 3, 4}]
>>> subsets(N)
{'11': set([2]), '10': set([0, 1]), '01': set([3, 4])}
>>> N = [{0, 1}, {2, 3}, {1, 3, 4}]
>>> subsets(N)
{'010': set([2]), '011': set([3]), '001': set([4]), '111': set([
]), '110': set([1]), '100': set([0]), '101': set([1])}
```

## 4.6 data\_tools.strings

String operations module.

data\_tools.strings.**is\_numeric**(s)

Determines if a string can be considered a numeric value. NaN is also considered, since it is float type.

- **Arguments:**

- s [str]: String to be evaluated.

- **Returns:**

- [bool]: True/False depending if the condition is satisfied.

- **Examples:**

```
>>> is_numeric('4')
True
>>> is_numeric('-3.2')
True
>>> is_numeric('number')
False
>>> is_numeric('NaN')
True
```

data\_tools.strings.**join\_str\_lists**(a, b, sep="")

Joins element-wise two lists (or any 1D iterable) of strings with a given separator (if provided). Length of the input lists must be equal.

- **Arguments:**

- a [list]: Contains the first elements [str] of the joint strings.
- b [list]: Contains the second elements [str] of the joint strings.
- sep [str]: Optional ' ' (non separated) by default. Determines the separator between the joint strings.

- **Returns:**

- [list]: List of the joint strings.

- **Example:**

```
>>> a = ['a', 'b']
>>> b = ['1', '2']
>>> join_str_lists(a, b, sep='_')
['a_1', 'b_2']
```





## PYTHON MODULE INDEX

### d

`data_tools.databases`, [9](#)  
`data_tools.diffusion`, [11](#)  
`data_tools.models`, [11](#)  
`data_tools.plots`, [13](#)  
`data_tools.sets`, [17](#)  
`data_tools.strings`, [19](#)



## B

bit\_or() (in module data\_tools.sets), 17

## C

chunk\_this() (in module data\_tools.sets), 17

## D

data\_tools.databases (module), 9  
 data\_tools.diffusion (module), 11  
 data\_tools.models (module), 11  
 data\_tools.plots (module), 13  
 data\_tools.sets (module), 17  
 data\_tools.strings (module), 19  
 density() (in module data\_tools.plots), 13  
 DoseResponse (class in data\_tools.models), 11

## E

ec() (data\_tools.models.DoseResponse method), 12  
 euler\_explicit1D() (in module data\_tools.diffusion), 11

## F

find\_min() (in module data\_tools.sets), 18  
 fit\_data() (data\_tools.models.Lasso method), 13

## I

in\_all() (in module data\_tools.sets), 18  
 is\_numeric() (in module data\_tools.strings), 19

## J

join\_str\_lists() (in module data\_tools.strings), 19

## K

kegg\_link() (in module data\_tools.databases), 9  
 kegg\_pathway\_mapping() (in module data\_tools.databases), 9

## L

Lasso (class in data\_tools.models), 12

## P

piano\_consensus() (in module data\_tools.plots), 14

plot() (data\_tools.models.DoseResponse method), 12  
 plot\_coef() (data\_tools.models.Lasso method), 13  
 plot\_score() (data\_tools.models.Lasso method), 13

## S

subsets() (in module data\_tools.sets), 18

## U

up\_map() (in module data\_tools.databases), 10

## V

venn() (in module data\_tools.plots), 15  
 volcano() (in module data\_tools.plots), 16