deepgraph Documentation

Release 0.2.3

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CHAPTER 1

Contents

1.1 What is DeepGraph

DeepGraph is an open source Python implementation of a new network representation introduced here. Its purpose is to facilitate data analysis by interpreting data in terms of network theory.

The basis of this software package is Pandas, a fast and flexible data analysis tool for the Python programming language. Utilizing one of its primary data structures, the DataFrame, we represent objects (i.e. the nodes of a network) by one DataFrame, and their pairwise relations (i.e. the edges of a network) by another DataFrame.

One of the main features of DeepGraph is an efficient and scalable creation of edges. Given a set of nodes in the form of a DataFrame (or an on disc HDFStore), DeepGraph's *core class* provides *methods* to iteratively compute pairwise relations between the nodes (e.g. similarity/distance measures) using arbitrary, user-defined functions on the nodes' features. These methods provide arguments to parallelize the computation and control memory consumption, making them suitable for very large data-sets and adjustable to whatever hardware you have at hand (from netbooks to cluster architectures).

Furthermore, once a graph is constructed, DeepGraph allows you to partition its *nodes*, *edges* or the entire *graph* by the graph's properties and labels, enabling the aggregation, computation and allocation of information on and between arbitrary *groups* of nodes. These methods also let you express elaborate queries on the information contained in a deep graph.

DeepGraph is not meant to replace or compete with already existing Python network libraries, such as NetworkX or graph_tool, but rather to combine and extend their capabilities with the merits of Pandas. For that matter, the core class of DeepGraph provides *interfacing methods* to convert to common network representations and graph objects of popular Python network packages.

Deepgraph also implements a number of useful plotting methods, including drawings on geographical map projections.

It's also possible to represent *multilayer networks* by deep graphs. We're thinking of implementing an interface to a suitable package dedicated to the analysis of multilayer networks.

Note: Please acknowledge the authors and cite the use of this software when results are used in publications or published elsewhere. Various citation formats are available here: https://aip.scitation.org/action/showCitFormats? type=show&doi=10.1063%2F1.4952963

For your convenience, you can find the BibTex entry below:

<pre>@Article{traxl-;</pre>	201	6-deep,
author	=	{Dominik Traxl AND Niklas Boers AND J\"urgen Kurths},
title	=	{Deep Graphs - A general framework to represent and analyze
		heterogeneous complex systems across scales},
journal	=	{Chaos},
year	=	{2016},
volume	=	{26},
number	=	{6},
eid	=	{065303},
doi	=	{http://dx.doi.org/10.1063/1.4952963},
eprinttype	=	{arxiv},
eprintclass	=	<pre>{physics.data-an, cs.SI, physics.ao-ph, physics.soc-ph},</pre>
eprint	=	{http://arxiv.org/abs/1604.00971v1},
version	=	{1},
date	=	{2016-04-04},
url	=	{http://arxiv.org/abs/1604.00971v1}
}		

To get started, have a look at

- Installation of DeepGraph
- DeepGraph's Tutorials
- API Reference

Want to share feedback, or contribute?

So far the package has only been developed by *me*, a fact that I would like to change very much. So if you feel like contributing in any way, shape or form, please feel free to contact me, report bugs, create pull requestes, milestones, etc. You can contact me via email: dominik.traxl@posteo.org

Note: This documentation assumes general familiarity with NumPy and Pandas. If you haven't used these packages, do invest some time in learning about them first.

Note: DeepGraph is free software; you can redistribute it and/or modify it under the terms of the BSD License. We highly welcome contributions from the community.

1.2 Installation

1.2.1 Quick Install

DeepGraph can be installed via pip from PyPI

```
$ pip install deepgraph
```

Depending on your system, you may need root privileges. On UNIX-based operating systems (Linux, Mac OS X etc.) this is achieved with sudo

\$ sudo pip install deepgraph

Alternatively, if you're using Conda, install with

\$ conda install -c conda-forge deepgraph

1.2.2 Installing from Source

Alternatively, you can install DeepGraph from source by downloading a source archive file (tar.gz or zip).

Source Archive File

- 1. Download the source (tar.gz or zip file) from https://pypi.python.org/pypi/deepgraph/ or https://github.com/ deepgraph/deepgraph/
- 2. Unpack and change directory to the source directory (it should have the files README.rst and setup.py).
- 3. Run python setup.py install to build and install. As a developer, you may want to install using cython: python setup.py install --use-cython.
- 4. (Optional) Run py.test to execute the tests if you have pytest installed.

GitHub

1. Clone the deepgraph repositiory

git clone https://github.com/deepgraph/deepgraph.git

- 2. Change directory to deepgraph
- 3. Run python setup.py install to build and install. As a developer, you may want to install using cython: python setup.py install --use-cython.
- 4. (Optional) Run py.test to execute the tests if you have pytest installed.

Installing without Root Privileges

If you don't have permission to install software on your system, you can install into another directory using the --user, --prefix, or --home flags to setup.py.

For example

\$ python setup.py install --prefix=/home/username/python

or

```
$ python setup.py install --home=~
```

or

\$ python setup.py install --user

Note: If you didn't install in the standard Python site-packages directory you will need to set your PYTHONPATH variable to the alternate location. See here for further details.

1.2.3 Requirements

The easiest way to get Python and the required/optional packages is to use Conda (or Miniconda), a cross-platform (Linux, Mac OS X, Windows) Python distribution for data analytics and scientific computing.

Python

To use DeepGraph you need Python 2.7, 3.4 or later.

Pandas

Pandas is an open source, BSD-licensed library providing high-performance, easy-to-use data structures and data analysis tools for the Python programming language.

Pandas is the core dependency of DeepGraph, and it is highly recommended to install the recommended and optional dependencies of Pandas as well.

NumPy

NumPy is the fundamental package for scientific computing with Python.

Needed for internal operations.

1.2.4 Recommended Packages

The following are recommended packages that DeepGraph can use to provide additional functionality.

Matplotlib

Matplotlib is a python 2D plotting library which produces publication quality figures in a variety of hardcopy formats and interactive environments across platforms.

Allows you to use the *plotting methods* of DeepGraph.

Matplotlib Basemap Toolkit

basemap is an add-on toolkit for matplotlib that lets you plot data on map projections with coastlines, lakes, rivers and political boundaries. See the basemap tutorial for documentation and examples of what it can do.

Used by plot_map and plot_map_generator to plot networks on map projections.

PyTables

PyTables is a package for managing hierarchical datasets and designed to efficiently and easily cope with extremely large amounts of data.

Necessary for HDF5-based storage of pandas DataFrames. DeepGraph's *core class* may be initialized with a HDFStore containing a node table in order to iteratively create edges directly from disc (see *create_edges* and *create_edges_ft*).

SciPy

SciPy is a Python-based ecosystem of open-source software for mathematics, science, and engineering.

Allows you to convert from DeepGraph's network representation to sparse adjacency matrices (see *return_cs_graph*).

NetworkX

NetworkX is a Python language software package for the creation, manipulation, and study of the structure, dynamics, and functions of complex networks.

Allows you to convert from DeepGraph's network representation to NetworkX's network representation (see *return_nx_graph*).

Graph-Tool

graph_tool is an efficient Python module for manipulation and statistical analysis of graphs (a.k.a. networks).

Allows you to convert from DeepGraph's network representation to Graph-Tool's network representation (see *return_gt_graph*).

Conda users can install graph_tool by adding the following channels to their ~/.condarc

```
$ conda config --add channels conda-forge
$ conda config --add channels ostrokach-forge
```

Then, install graph-tool

```
$ conda install graph-tool
```

You can test your graph-tool installation by

```
$ python -c "from graph_tool.all import *"
```

1.2.5 Optional Packages

The following packages are considered to provide very useful tools and methods.

Scikit-Learn

sklearn is a Python module integrating classical machine learning algorithms in the tightly-knit world of scientific Python packages (numpy, scipy, matplotlib).

Sklearn-pandas

sklearn-pandas provides a bridge between Scikit-Learn's machine learning methods and pandas-style Data Frames.

1.3 Tutorials

1.3.1 10 Minutes to DeepGraph

[ipython notebook] [python script] [data]

This is a short introduction to DeepGraph. In the following, we demonstrate DeepGraph's core functionalities by a toy data-set, "flying balls".

First of all, we need to import some packages

```
# for plots
import matplotlib.pyplot as plt
# the usual
import numpy as np
import pandas as pd
import deepgraph as dg
# notebook display
%matplotlib inline
plt.rcParams['figure.figsize'] = 8, 6
pd.options.display.max_rows = 10
pd.set_option('expand_frame_repr', False)
```

Loading Toy Data

Then, we need data in the form of a pandas DataFrame, representing the nodes of our graph

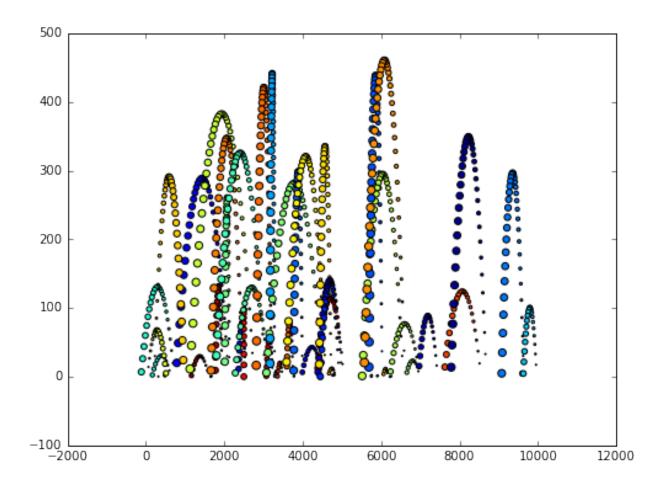
```
v = pd.read_csv('flying_balls.csv', index_col=0)
print(v)
```

	time	Х	У	ball_id
0	0	1692.000000	0.000000	0
1	0	8681.000000	0.00000	1
2	0	490.000000	0.00000	2
3	0	7439.000000	0.000000	3
4	0	4998.000000	0.00000	4
1163	45	2812.552734	16.503178	39
1164	46	5686.915998	14.161693	10
1165	46	3161.729086	19.381823	14
1166	46	5594.233413	57.701712	37
1167	47	5572.216748	20.588750	37
[1168	rows	x 4 columns]		

The data consists of 1168 space-time measurements of 50 different toy balls in two-dimensional space. Each space-time measurement (i.e. row of v) represents a **node**.

Let's plot the data such that each ball has it's own color

```
plt.scatter(v.x, v.y, s=v.time, c=v.ball_id)
```



Creating Edges

In order to create edges between these nodes, we now initiate a dg. DeepGraph instance

```
g = dg.DeepGraph(v)
g
```

<DeepGraph object, with n=1168 node(s) and m=0 edge(s) at 0x7facf3b35dd8>

and use it to create edges between the nodes given by q. v. For that matter, we may define a **connector** function

```
def x_dist(x_s, x_t):
    dx = x_t - x_s
    return dx
```

and pass it to g.create_edges in order to compute the distance in the x-coordinate of each pair of nodes

```
g.create_edges(connectors=x_dist)
g
```

<DeepGraph object, with n=1168 node(s) and m=681528 edge(s) at 0x7facf3b35dd8>

print(g.e)

		dx
s	t	
0	1	6989.000000
	2	-1202.000000
	3	5747.000000
	4	3306.000000
	5	2812.000000
1164	1166	-92.682585
	1167	-114.699250
1165	1166	2432.504327
	1167	2410.487662
1166	1167	-22.016665
[6815	528 ro	ows x 1 columns]

Let's say we're only interested in creating edges between nodes with a x-distance smaller than 1000. Then we may additionally define a **selector**

```
def x_dist_selector(dx, sources, targets):
    dxa = np.abs(dx)
    sources = sources[dxa <= 1000]
    targets = targets[dxa <= 1000]
    return sources, targets</pre>
```

and pass both the **connector** and **selector** to *g.create_edges*

g.create_edges(connectors=x_dist, selectors=x_dist_selector)

<DeepGraph object, with n=1168 node(s) and m=156938 edge(s) at 0x7facf3b35dd8>

print(g.e)

g

		dx
s	t	
0	6	416.000000
	7	848.000000
	19	-973.000000
	24	437.000000
	38	778.000000
1162	1167	-44.033330
1163	1165	349.176351
1164	1166	-92.682585
	1167	-114.699250
1166	1167	-22.016665
[156	938 ro	ows x 1 columns]

There is, however, a much more efficient way of creating edges that involve a simple distance threshold such as the one above

Creating Edges on a FastTrack

In order to efficiently create edges including a selection of edges via a simple distance threshold as above, one should use the *create_edges_ft* method. It relies on a sorted DataFrame, so we need to sort *g*. v first

```
g.v.sort_values('x', inplace=True)
```

```
g.create_edges_ft(ft_feature=('x', 1000))
g
```

<DeepGraph object, with n=1168 node(s) and m=156938 edge(s) at 0x7facf3b35dd8>

Let's compare the efficiency

%timeit -n3 -r3 g.create_edges(connectors=x_dist, selectors=x_dist_selector)

3 loops, best of 3: 557 ms per loop

%timeit -n3 -r3 g.create_edges_ft(ft_feature=('x', 1000))

3 loops, best of 3: 167 ms per loop

The *create_edges_ft* method also accepts **connectors** and **selectors** as input. Let's connect only those measurements that are close in space and time

```
def y_dist(y_s, y_t):
    dy = y_t - y_s
    return dy
def time_dist(time_t, time_s):
    dt = time_t - time_s
    return dt
def y_dist_selector(dy, sources, targets):
    dya = np.abs(dy)
    sources = sources[dya <= 100]</pre>
    targets = targets[dya <= 100]</pre>
    return sources, targets
def time_dist_selector(dt, sources, targets):
   dta = np.abs(dt)
    sources = sources[dta <= 1]</pre>
    targets = targets[dta <= 1]</pre>
    return sources, targets
```

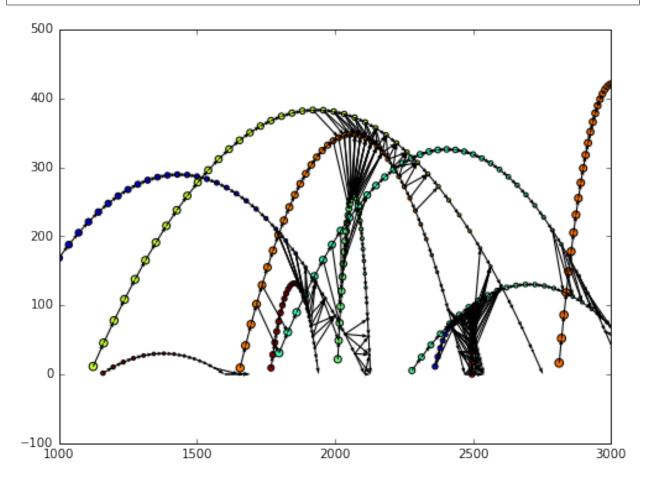
g

<DeepGraph object, with n=1168 node(s) and m=1899 edge(s) at 0x7facf3b35dd8>

print(g.e)

		dt	dy	ft_r
S	t			
890	867	-1	19.311136	33.415831
867	843	-1	17.678482	33.415831
843	818	-1	16.045829	33.415831
818	792	-1	14.413176	33.415831
792	766	-1	12.780523	33.415831
		• •		
244	203	-1	-10.825226	15.455612
203	159	-1	-12.457879	15.455612
159	114	-1	-14.090532	15.455612
114	65	-1	-15.723185	15.455612
65	16	-1	-17.355838	15.455612
[189	99 ro	WS >	k 3 columns]	

We can now plot the flying balls and the edges we just created with the plot_2d method



Graph Partitioning

The *DeepGraph* class also offers methods to partition *nodes*, *edges* and an entire *graph*. See the docstrings and the other tutorials for details and examples.

Graph Interfaces

Furthermore, you may inspect the docstrings of *return_cs_graph*, *return_nx_graph* and *return_gt_graph* to see how to convert from DeepGraph's DataFrame representation of a network to sparse adjacency matrices, NetworkX's network representation and graph_tool's network representation.

Plotting Methods

DeepGraph also offers a number of useful Plotting methods. See *plotting methods* for details and have a look at the other tutorials for examples.

1.3.2 Computing Very Large Correlation Matrices in Parallel

[ipython notebook] [python script]

Note: Please acknowledge the authors and cite the use of this software when results are used in publications or published elsewhere. Various citation formats are available here: https://aip.scitation.org/action/showCitFormats? type=show&doi=10.1063%2F1.4952963

For your convenience, you can find the BibTex entry below:

```
@Article{trax1-2016-deep,
```

```
author = {Dominik Traxl AND Niklas Boers AND J\"urgen Kurths},
           = {Deep Graphs - A general framework to represent and analyze
title
               heterogeneous complex systems across scales},
journal
          = \{Chaos\},
           = \{2016\},\
year
volume
           = \{26\},\
number
            = \{6\},\
eid
            = \{065303\},\
           = {http://dx.doi.org/10.1063/1.4952963},
doi
eprinttype = {arxiv},
eprintclass = {physics.data-an, cs.SI, physics.ao-ph, physics.soc-ph},
eprint = {http://arxiv.org/abs/1604.00971v1},
version
           = \{1\},
           = \{2016 - 04 - 04\},\
date
            = {http://arxiv.org/abs/1604.00971v1}
url
```

In this short tutorial, we'll demonstrate how DeepGraph can be used to efficiently compute very large correlation matrices in parallel, with full control over RAM usage.

Assume you have a set of n_samples samples, each comprised of n_features features and you want to compute the Pearson correlation coefficients between all pairs of features (for the Spearman's rank correlation coefficients, see the *Note*-box below). If your data is small enough, you may use scipy.stats.pearsonr or numpy.corrcoef, but for large data, neither of these methods is feasible. Scipy's pearsonr would be very slow, since you'd have to compute pair-wise correlations in a double loop, and numpy's corrcoef would most likely blow your RAM.

Using DeepGraph's *create_edges* method, you can compute all pair-wise correlations efficiently. In this tutorial, the data is stored on disc and only the relevant subset of features for each iteration will be loaded into memory by the computing nodes. Parallelization is achieved by using python's standard library multiprocessing, but it should be straight-forward to modify the code to accommodate other parallelization libraries. It should also be straight-forward to modify the code in order to compute other correlation/distance/similarity-measures between a set of features.

First of all, we need to import some packages

```
# data i/o
import os
# compute in parallel
from multiprocessing import Pool
# the usual
import numpy as np
import pandas as pd
import deepgraph as dg
```

Let's create a set of variables and store it as a 2d-matrix X (shape=(n_features, n_samples)) on disc. To speed up the computation of the correlation coefficients later on, we whiten each variable.

Note: On the computation of the Spearman's rank correlation coefficients: Since the Spearman correlation coefficient is defined as the Pearson correlation coefficient between the ranked variables, it suffices to uncomment the indicated line in the above code-block in order to compute the Spearman's rank correlation coefficients in the following.

Now we can compute the pair-wise correlations using DeepGraph's *create_edges* method. Note that the node table v only stores references to the mem-mapped array containing the samples.

```
# parameters (change these to control RAM usage)
step_size = 1e5
n_processes = 100
# load samples as memory-map
X = np.load('samples.npy', mmap_mode='r')
# create node table that stores references to the mem-mapped samples
v = pd.DataFrame({'index': range(X.shape[0])})
```

```
(continued from previous page)
```

```
# connector function to compute pairwise pearson correlations
def corr(index_s, index_t):
   features_s = X[index_s]
   features_t = X[index_t]
   corr = np.einsum('ij,ij->i', features_s, features_t) / n_samples
   return corr
# index array for parallelization
pos_array = np.array(np.linspace(0, n_features*(n_features-1)//2, n_processes),_
→dtype=int)
# parallel computation
def create_ei(i):
   from_pos = pos_array[i]
   to_pos = pos_array[i+1]
    # initiate DeepGraph
   q = dq.DeepGraph(v)
    # create edges
   g.create_edges(connectors=corr, step_size=step_size,
                   from_pos=from_pos, to_pos=to_pos)
    # store edge table
   g.e.to_pickle('tmp/correlations/{}.pickle'.format(str(i).zfill(3)))
# computation
if __name__ == '__main__':
   os.makedirs("tmp/correlations", exist_ok=True)
   indices = np.arange(0, n_processes - 1)
   p = Pool()
   for _ in p.imap_unordered(create_ei, indices):
       pass
```

Let's collect the computed correlation values and store them in an hdf file.

```
# store correlation values
files = os.listdir('tmp/correlations/')
files.sort()
store = pd.HDFStore('e.h5', mode='w')
for f in files:
    et = pd.read_pickle('tmp/correlations/{}'.format(f))
    store.append('e', et, format='t', data_columns=True, index=False)
store.close()
```

Let's have a quick look at the correlations.

```
# load correlation table
e = pd.read_hdf('e.h5')
print(e)
```

```
corr
s t
0 1 -0.006066
```

	2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29	0.094063 -0.025529 0.074080 0.035490 0.005221 0.032064 0.000378 -0.049318 -0.084853 0.026407 0.028543 -0.013347 -0.180113 0.151164 -0.094398 -0.124582 -0.000781 -0.044138 -0.193609 0.003877 0.048305 0.004877 -0.021291 -0.070756 -0.014906 -0.197605 -0.103509 0.071503
 4991 4992	30 4998 4999 4993 4994 4995	0.120718
4993	4996 4997 4998 4999 4994 4995 4996 4997	-0.052319 -0.010797 -0.025321
4994	4998 4999 4995 4996 4997	-0.069946 -0.031208 0.108063 0.144441 0.078353
4995	4998 4999 4996 4997 4998	-0.024799 -0.026432 -0.019991 -0.178458 -0.162406
4996	4999 4997 4998 4999	0.102835 0.115812 -0.061167 0.018606

```
4997 4998 -0.151932
4999 -0.271358
4998 4999 0.106453
[12497500 rows x 1 columns]
```

And finally, let's see where most of the computation time is spent.

```
g = dg.DeepGraph(v)
p = %prun -r g.create_edges(connectors=corr, step_size=step_size)
```

p.print_stats(20)

```
244867 function calls (239629 primitive calls) in 14.193 seconds
  Ordered by: internal time
  List reduced from 541 to 20 due to restriction <20>
  ncalls tottime percall cumtime percall filename:lineno(function)
     250
           9.355
                   0.037
                            9.361
                                      0.037 memmap.py:334 (__getitem__)
     125
            1.584
                    0.013
                             1.584
                                      0.013 {built-in method numpy.core.multiarray.
\hookrightarrow c einsum}
     125
            1.012
                  0.008 12.013
                                      0.096 deepgraph.py:4558(map)
            0.581
                  0.290
                           0.581
       2
                                     0.290 {method 'get_labels' of 'pandas._libs.
→hashtable.Int64HashTable' objects}
          0.301 0.301 0.414
       1
                                      0.414 multi.py:795(_engine)
       5
           0.157 0.031
                             0.157 0.031 {built-in method numpy.core.multiarray.
\hookrightarrow concatenate \}
     250
           0.157
                  0.001 0.170
                                      0.001 internals.py:5017(_stack_arrays)
           0.105
                  0.053
                            0.105
       2.
                                      0.053 {pandas._libs.algos.take_1d_int64_int64}
                                      0.000 {method 'reduce' of 'numpy.ufunc'
           0.094
                    0.000
                            0.094
     889
→objects}
     125
            0.089
                    0.001 12.489
                                      0.100 deepgraph.py:5294 (_select_and_return)
     125
           0.074
                    0.001
                             0.074
                                      0.001 {deepgraph._triu_indices._reduce_triu_
\rightarrow indices}
     125
           0.066
                    0.001
                             0.066
                                     0.001 {built-in method deepgraph._triu_

→indices._triu_indices}

                           0.038
       4
           0.038 0.009
                                    0.009 {built-in method pandas._libs.algos.
\rightarrowensure_int16}
     125
          0.033 0.000
                           10.979
                                    0.088 <ipython-input-3-26c4f59cd911>:12(corr)
           0.028 0.014 0.028 0.014 function_base.py:4703(delete)
       2
       1
           0.027 0.027 14.163 14.163 deepgraph.py:4788(_matrix_iterator)
            0.027
                    0.027 0.113 0.113 multi.py:56(_codes_to_ints)
       1
45771/45222
            0.020
                    0.000 0.043
                                        0.000 {built-in method builtins.isinstance}
       1
            0.019
                    0.019 14.193 14.193 deepgraph.py:183(create_edges)
       2
            0.012
                    0.006
                             0.700
                                      0.350 algorithms.py:576(factorize)
```

As you can see, most of the time is spent by getting the requested features in the corr-function, followed by computing the correlation values themselves.

1.3.3 Building a DeepGraph of Extreme Precipitation

[ipython notebook] [python script] [data]

In the following we build a deep graph of a high-resolution dataset of precipitation measurements.

The goal is to first detect spatiotemporal clusters of extreme precipitation events and then to create families of these clusters based on a spatial correlation measure. Finally, we create and plot some informative (intersection) partitions of the deep graph.

For further details see section V of the original paper: https://arxiv.org/abs/1604.00971

First of all, we need to import some packages

```
# data i/o
import os
import os
import xarray
# for plots
import matplotlib.pyplot as plt
# the usual
import numpy as np
import pandas as pd
import deepgraph as dg
# notebook display
from IPython.display import HTML
%matplotlib inline
plt.rcParams['figure.figsize'] = 8, 6
pd.options.display.max_rows = 10
pd.set_option('expand_frame_repr', False)
```

Selecting and Preprocessing the Precipitation Data

Selection

If you want to select your own spatiotemporal box of precipitation events, you may follow the instructions below and change the filename in the next box of code.

- Go to https://disc.gsfc.nasa.gov/datasets/TRMM_3B42_V7/summary?keywords=TRMM_3B42_V7
- click on "Simple Subset Wizard"
- select the "Date Range" (and if desired a "Spatial Bounding Box") you're interested in
- click on "Search for Data Sets"
- expand the list by clicking on the "+" symbol
- mark the check box "precipitation"
- (optional, but recommended) click on the selector to change from "netCDF" to "gzipped netCDF"
- click on "Subset Selected Data Sets"
- click on "View Subset Results"
- right click on the "Get list of URLs for this subset in a file" link, and choose "Save Link As..."
- the downloaded file will have a name similar to "SSW_download_2016-05-03T20_19_28_23621_20Ie06xp.inp". Note which directory the downloaded file is saved to, and in your Unix shell, set your current working directory to that directory.
- register an account to get authentication credentials using these instructions: https://disc.gsfc.nasa.gov/ information/howto/5761bc6a5ad5a18811681bae?keywords=wget

· get the files via

```
os.system("wget --content-disposition --directory-prefix=tmp --load-cookies ~/.urs_

→cookies --save-cookies ~/.urs_cookies --auth-no-challenge=on --keep-session-cookies_

→-i SSW_download_2016-05-03T20_19_28_23621_20Ie06xp.inp")
```

Preprocessing

Next, we need to convert the downloaded netCDF files to a pandas DataFrame, which we can then use to initiate a *dg.DeepGraph*

```
# choose "wet times" threshold
r = .1
# choose "extreme" precipitation threshold
p = .9
v_list = []
for file in os.listdir('tmp'):
   if file.startswith('3B42.'):
        # open the downloaded netCDF file
        # unfortunately, we have to decode times ourselves, since
        # the format of the downloaded files doesn't work
        # see also: https://github.com/pydata/xarray/issues/521
        f = xarray.open_dataset('tmp/{}'.format(file), decode_times=False)
        # create integer-based (x,y) coordinates
        f['x'] = (('longitude'), np.arange(len(f.longitude)))
        f['y'] = (('latitude'), np.arange(len(f.latitude)))
        # convert to pd.DataFrame
        vt = f.to_dataframe()
        # we only consider "wet times", pcp >= 0.1mm/h
       vt = vt[vt.pcp >= r]
        # reset index
        vt.reset_index(inplace=True)
        # add correct times
        ftime = f.time.units.split()[2:]
        year, month, day = ftime[0].split('-')
        hour = ftime[1]
        time = pd.datetime(int(year), int(month), int(day), int(hour))
        vt['time'] = time
        # compute "area" for each event
        vt['area'] = 111**2 * .25**2 * np.cos(2*np.pi*vt.latitude / 360.)
        # compute "volume of water precipitated" for each event
        vt['vol'] = vt.pcp * 3 * vt.area
        # set dtypes -> economize ram
        vt['pcp'] = vt['pcp'] * 100
        vt['pcp'] = vt['pcp'].astype(np.uint16)
        vt['latitude'] = vt['latitude'].astype(np.float16)
```

```
(continued from previous page)
```

```
vt['longitude'] = vt['longitude'].astype(np.float16)
        vt['area'] = vt['area'].astype(np.uint16)
        vt['vol'] = vt['vol'].astype(np.uint32)
        vt['x'] = vt['x'].astype(np.uint16)
        vt['y'] = vt['y'].astype(np.uint16)
        # append to list
        v_list.append(vt)
        f.close()
# concatenate the DataFrames
v = pd.concat(v_list)
# append a column indicating geographical locations (i.e., supernode labels)
v['g_id'] = v.groupby(['longitude', 'latitude']).grouper.group_info[0]
v['g_id'] = v['g_id'].astype(np.uint32)
# select `p`th percentile of precipitation events for each geographical location
v = v.groupby('g_id').apply(lambda x: x[x.pcp >= x.pcp.quantile(p)])
# append integer-based time
dtimes = pd.date_range(v.time.min(), v.time.max(), freq='3H')
dtdic = {dtime: itime for itime, dtime in enumerate(dtimes)}
v['itime'] = v.time.apply(lambda x: dtdic[x])
v['itime'] = v['itime'].astype(np.uint16)
# sort by time
v.sort_values('time', inplace=True)
# set unique node index
v.set_index(np.arange(len(v)), inplace=True)
# shorten column names
v.rename(columns={'pcp': 'r',
                  'latitude': 'lat',
                  'longitude': 'lon',
                  'time': 'dtime',
                  'itime': 'time'},
         inplace=True)
```

The created DataFrame of extreme precipitation measurements looks like this

	lat	lon	dtime	r	Х	У	area	vol	g_id	time
0	15.125	-118.125	2004-08-20	1084	28	101	743	24174	5652	0
1	44.875	-30.625	2004-08-20	392	378	220	545	6433	85341	0
2	45.125	-30.625	2004-08-20	454	378	221	543	7416	85342	0
3	45.375	-30.625	2004-08-20	909	378	222	540	14767	85343	0
4	45.625	-30.625	2004-08-20	907	378	223	538	14669	85344	0
382306	26.875	-46.625	2004-09-27	503	314	148	686	10385	70380	304
382307	38.375	-37.125	2004-09-27	453	352	194	603	8222	79095	304
382308	8.125	-105.125	2004-09-27	509	80	73	762	11663	17007	304
382309	21.875	-42.875	2004-09-27	260	329	128	714	5595	73875	304
382310	6.625	-111.125	2004-09-27	192	56	67	764	4428	11790	304
										(continues on next nage

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print(v)

[382311 rows x 10 columns]

We identify each row of this table as a node of our *DeepGraph*

```
g = dg.DeepGraph(v)
```

Plot the Data

Let's take a look at the data by creating a video of the time-evolution of precipitation measurements. Using the *plot_map_generator* method, this is straight forward.

```
# configure map projection
kwds_basemap = {'llcrnrlon': v.lon.min() - 1,
                'urcrnrlon': v.lon.max() + 1,
                'llcrnrlat': v.lat.min() - 1,
                'urcrnrlat': v.lat.max() + 1,
                'resolution': 'i'}
# configure scatter plots
kwds_scatter = {'s': 1.5,
                'c': g.v.r.values / 100.,
                'edgecolors': 'none',
                'cmap': 'viridis_r'}
# create generator of scatter plots on map
objs = g.plot_map_generator('lon', 'lat', 'dtime',
                            kwds_basemap=kwds_basemap,
                            kwds_scatter=kwds_scatter)
# plot and store frames
for i, obj in enumerate(objs):
    # configure plots
   cb = obj['fig'].colorbar(obj['pc'], fraction=0.025, pad=0.01)
   cb.set_label('[mm/h]')
   obj['m'].fillcontinents(color='0.2', zorder=0, alpha=.4)
   obj['ax'].set_title('{}'.format(obj['group']))
    # store and close
   obj['fig'].savefig('tmp/pcp_{:03d}.png'.format(i),
                       dpi=300, bbox_inches='tight')
    plt.close(obj['fig'])
```

<	/	v	ideo>
n	"	")

[download video]

Detecting SpatioTemporal Clusters of Extreme Precipitation

In this tutorial, we're interested in local formations of spatiotemporal clusters of extreme precipitation events. For that matter, we now use DeepGraph to identify such clusters and track their temporal evolution.

Create Edges

We now use DeepGraph to create edges between the nodes given by g.v.

The edges of g will be utilized to detect spatiotemporal clusters in the data, or in more technical terms: to partition the set of all nodes into subsets of connected grid points. One can imagine the nodes to be elements of a 3 dimensional grid box (x,y,time), where we allow every node to have 26 possible neighbours (8 neighbours in the time slice of the measurement, t_i , and 9 neighbours in each the time slice t_i 1 and $t_i + 1$).

For that matter, we pass the following connectors

```
def grid_2d_dx(x_s, x_t):
    dx = x_t - x_s
    return dx
def grid_2d_dy(y_s, y_t):
    dy = y_t - y_s
    return dy
```

and selectors

```
def s_grid_2d_dx(dx, sources, targets):
    dxa = np.abs(dx)
    sources = sources[dxa <= 1]
    targets = targets[dxa <= 1]
    return sources, targets

def s_grid_2d_dy(dy, sources, targets):
    dya = np.abs(dy)
    sources = sources[dya <= 1]
    targets = targets[dya <= 1]
    return sources, targets</pre>
```

to the <code>create_edges_ft</code> method

```
# rename fast track relation
g.e.rename(columns={'ft_r': 'dt'}, inplace=True)
```

To see how many nodes and edges our graph's comprised of, one may simply type

g

```
<DeepGraph object, with n=382311 node(s) and m=567225 edge(s) at 0x7f7a4c3de160>
```

The edges we just created look like this

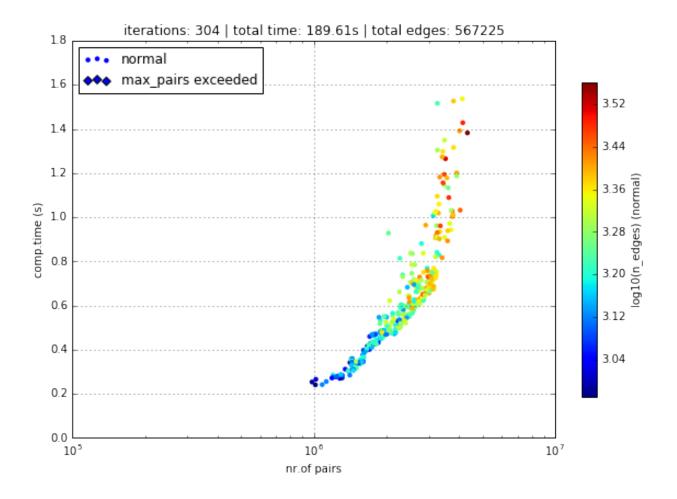
```
print(g.e)
```

		dx	dy	dt
S	t			
0	1362	0	1	False
	1432	1	0	False
	1433	1	1	False
	1696	1	0	True
	1699	1	1	True
382284	382291	0	1	False
382295	382296	0	1	False
382296	382299	0	1	False
382299	382309	0	1	False
382304	382306	0	1	False
[56722	5 rows x	3 c	olum	ins]

Logfile Plot

To see how long it took to create the edges, one may use the *plot_logfile* method

g.plot_logfile('create_e')



Find the Connected Components

Having linked all neighbouring nodes, the spatiotemporal clusters can be identified as the connected components of the graph. For practical reasons, *DeepGraph* directly implements a method to find the connected components of a graph, *append_cp*

```
# all singular components (components comprised of one node only)
# are consolidated under the label 0
g.append_cp(consolidate_singles=True)
# we don't need the edges any more
del g.e
```

the node table now has a component membership column appended

print(print(g.v)													
	lat	lon	dtime	r	x	У	area	vol	g_id	time	ср			
0	15.125	-118.125	2004-08-20	1084	28	101	743	24174	5652	0	865			
1	44.875	-30.625	2004-08-20	392	378	220	545	6433	85341	0	5079			
2	45.125	-30.625	2004-08-20	454	378	221	543	7416	85342	0	5079			
3	45.375	-30.625	2004-08-20	909	378	222	540	14767	85343	0	5079			
4	45.625	-30.625	2004-08-20	907	378	223	538	14669	85344	0	5079			
•••	• • •	• • •		• • •	• • •	• • •	• • •			• • •				

									(cont	inued fro	m previous page)
382306	26.875	-46.625	2004-09-27	503	314	148	686	10385	70380	304	609
382307	38.375	-37.125	2004-09-27	453	352	194	603	8222	79095	304	0
382308	8.125	-105.125	2004-09-27	509	80	73	762	11663	17007	304	174
382309	21.875	-42.875	2004-09-27	260	329	128	714	5595	73875	304	8
382310	6.625	-111.125	2004-09-27	192	56	67	764	4428	11790	304	15610
[382311	rows x	11 column	ıs]								

Let's see how many spatiotemporal clusters g is comprised of (discarding singular components)

g.v.cp.max()

and how many nodes there are in the components

```
print(g.v.cp.value_counts())
```

Partition the Nodes Into a Component Supernode Table

In order to aggregate and compute some information about the precipitiation clusters, we now partition the nodes by the type of feature cp, the component membership labels of the nodes just created. This can be done with the *partition_nodes* method

```
# feature functions, will be applied to each component of g
feature_funcs = {'dtime': [np.min, np.max],
                          'time': [np.min, np.max],
                         'vol': [np.sum],
                          'lat': [np.mean],
                          'lon': [np.mean]}
# partition the node table
cpv, gv = g.partition_nodes('cp', feature_funcs, return_gv=True)
# append geographical id sets
cpv['g_ids'] = gv['g_id'].apply(set)
# append cardinality of g_id sets
cpv['n_unique_g_ids'] = cpv['g_ids'].apply(len)
# append time spans
```

```
cpv['dt'] = cpv['dtime_amax'] - cpv['dtime_amin']
# append spatial coverage
def area(group):
    return group.drop_duplicates('g_id').area.sum()
cpv['area'] = gv.apply(area)
```

The clusters look like this

print(cpv)

n_nodes dtime_amin dtime_amax time_amin time_	amax lat_	
→mean vol_sum lon_mean	g_ids n_	
→unique_g_ids dt area		
ср		
0 64678 2004-08-20 00:00:00 2004-09-27 00:00:00 0		
↔609375 627097323 -63.40625 {0, 1, 2, 6, 7, 10, 12, 13, 14, 22, 23,	24, 25	<u>ل</u>
↔ 49808 38 days 00:00:00 34781178		
1 16460 2004-09-01 06:00:00 2004-09-17 18:00:00 98		
→281250 351187150 -65.12500 {65536, 65537, 65538, 65539, 65540, 6554	1, 655	<u>ں</u>
→ 6629 16 days 12:00:00 4803624		
2 8519 2004-09-17 03:00:00 2004-09-24 15:00:00 225		
→906250 133698579 -44.62500 {73728, 73729, 73730, 73731, 73732, 7373	3, 737	_
→ 3730 7 days 12:00:00 2507350		
3 6381 2004-08-26 09:00:00 2004-09-06 03:00:00 51		
↔062500 113782748 -64.12500 {65555, 65556, 65557, 65558, 65559, 6556	0, 655	ш
→ 2442 10 days 18:00:00 1749673		
4 3403 2004-08-21 21:00:00 2004-08-24 12:00:00 15		
↔578125 66675326 -111.93750 {8141, 14654, 11805, 16363, 8142, 11806,	20490	_
→ 1294 2 days 15:00:00 978604		
•••• ••• ••• •••	•••	•
↔	· · · ·	
···· ···		
33165 2 2004-08-23 18:00:00 2004-08-23 18:00:00 30		
↔500000 20212 -103.87500 {1811	5, 18116} <u></u>	L
→ 2 0 days 00:00:00 1483		
33166 2 2004-09-05 18:00:00 2004-09-05 18:00:00 134		
↔250000 9366 -121.87500 {26	88, 268/}	ш
→ 2 0 days 00:00:00 1368	0.5	
33167 2 2004-08-30 15:00:00 2004-08-30 15:00:00 85		
$\Rightarrow 250000 \qquad 43096 \qquad 0.62500 \qquad \{112116$, 11211/}	—
→ 2 0 days 00:00:00 1519	1.01 0	
33168 2 2004-09-09 03:00:00 2004-09-09 03:00:00 161 →750000 24156 -13.62500 {100613		
	, 100614}	-
→ 2 0 days 00:00:00 1528 33169 2 2004-09-11 03:00:00 2004-09-11 03:00:00 177	177 15	
	2 005241	
	J, 90024}	-
→ 2 0 days 00:00:00 1483		
[33170 rows x 12 columns]		Ĩ

Plot the Largest Component

Let's see how the largest cluster of extreme precipitation evolves over time, again using the *plot_map_generator* method

```
# temporary DeepGraph instance containing
# only the largest component
gt = dg.DeepGraph(g.v)
gt.filter_by_values_v('cp', 1)
# configure map projection
from mpl_toolkits.basemap import Basemap
m1 = Basemap(projection='ortho',
             lon_0=cpv.loc[1].lon_mean + 12,
             lat_0=cpv.loc[1].lat_mean + 8,
             resolution=None)
width = (ml.urcrnrx - ml.llcrnrx) * .65
height = (m1.urcrnry - m1.llcrnry) * .45
kwds_basemap = {'projection': 'ortho',
                'lon_0': cpv.loc[1].lon_mean + 12,
                'lat_0': cpv.loc[1].lat_mean + 8,
                'llcrnrx': -0.5 * width,
                'llcrnry': -0.5 * height,
                'urcrnrx': 0.5 * width,
                'urcrnry': 0.5 * height,
                'resolution': 'i'}
# configure scatter plots
kwds_scatter = {'s': 2,
                'c': np.log(gt.v.r.values / 100.),
                'edgecolors': 'none',
                'cmap': 'viridis_r'}
# create generator of scatter plots on map
objs = gt.plot_map_generator('lon', 'lat', 'dtime',
                              kwds_basemap=kwds_basemap,
                              kwds_scatter=kwds_scatter)
# plot and store frames
for i, obj in enumerate(objs):
    # configure plots
   obj['m'].fillcontinents(color='0.2', zorder=0, alpha=.4)
   obj['m'].drawparallels(range(-50, 50, 20), linewidth=.2)
   obj['m'].drawmeridians(range(0, 360, 20), linewidth=.2)
   obj['ax'].set_title('{}'.format(obj['group']))
    # store and close
   obj['fig'].savefig('tmp/cp1_ortho_{:03d}.png'.format(i),
                       dpi=300, bbox_inches='tight')
   plt.close(obj['fig'])
```

```
</video>
""")
```

[download video]

Detecting Families of Spatially Related Clusters

Create SuperEdges between the Components

We now create superedges between the spatiotemporal clusters in order to find families of clusters that have a strong regional overlap. Passing the following **connectors** and **selector**

```
# compute intersection of geographical locations
def cp_node_intersection(g_ids_s, g_ids_t):
   intsec = np.zeros(len(g_ids_s), dtype=object)
   intsec_card = np.zeros(len(q_ids_s), dtype=np.int)
   for i in range(len(g ids s)):
        intsec[i] = g_ids_s[i].intersection(g_ids_t[i])
        intsec_card[i] = len(intsec[i])
    return intsec_card
# compute a spatial overlap measure between clusters
def cp intersection strength (n unique q ids s, n unique q ids t, intsec card):
   min_card = np.array(np.vstack((n_unique_g_ids_s, n_unique_g_ids_t)).min(axis=0),
                        dtype=np.float64)
    intsec_strength = intsec_card / min_card
   return intsec_strength
# compute temporal distance between clusters
def time_dist(dtime_amin_s, dtime_amin_t):
   dt = dtime_amin_t - dtime_amin_s
   return dt
```

to the *create_edges* method will provide the information necessary for this task

Since no selection of edges has taken place, the number of edges should be cpg.n*(cpg.n-1)/2

cpg

<DeepGraph object, with n=5000 node(s) and m=12497500 edge(s) at 0x7f7a00aec128>

print(cpg.e)

		intsec_strength
s	t	
1	2	0.018499
	3	0.002457
	4	0.00000
	5	0.00000
	6	0.00000
4997	4999	0.00000
	5000	0.00000
4998	4999	0.00000
	5000	0.00000
4999	5000	0.00000
[1249	97500	rows x 1 columns]

print(cpg.e.intsec_strength.value_counts())

0.000000	12481941		
1.000000	787		
0.111111	488		
0.333333	481		
0.500000	462		
0.012346	1		
0.158537	1		
0.178082	1		
0.658537	1		
0.018809	1		
Name: int	sec_strength,	dtype:	int64

Hierarchically Agglomerate Clusters into Families

Based on the above measure of spatial overlap between clusters, we now perform an agglomerative, hierarchical clustering of the spatio-temporal clusters into regionally coherent families.

```
from scipy.cluster.hierarchy import linkage, fcluster
# create condensed distance matrix
dv = 1 - cpg.e.intsec_strength.values
del cpg.e
# create linkage matrix
lm = linkage(dv, method='average', metric='euclidean')
del dv
# form flat clusters and append their labels to cpv
cpv['F'] = fcluster(lm, 1000, criterion='maxclust')
del lm
```

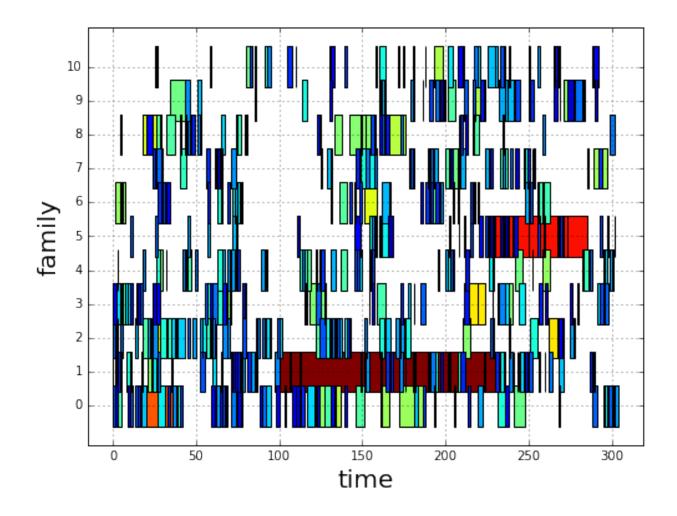
```
# relabel families by size
f = cpv['F'].value_counts().index.values
fdic = {j: i for i, j in enumerate(f)}
cpv['F'] = cpv['F'].apply(lambda x: fdic[x])
```

Let's see how many clusters there are in the families

```
print(cpv['F'].value_counts())
```

Create a "Raster Plot" of Families

Let's plot the clusters of the largest 10 families in a raster-like boxplot, by means of the plot_rects_label_numeric method



Create and Plot Informative (Intersection) Partitions

In this last section, we create some useful (intersection) partitions of the deep graph, which we then use to create some plots.

Geographical Locations

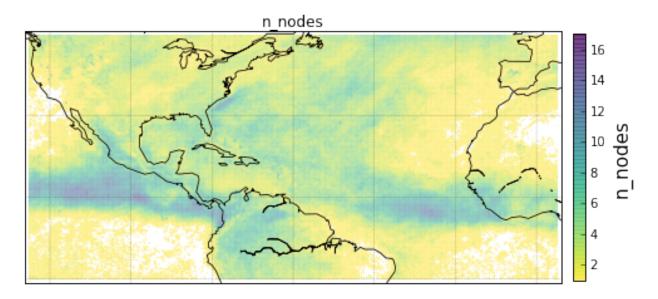
```
# how many components have hit a certain
# geographical location (discarding singular cps)
def count(cp):
    return len(set(cp[cp != 0]))
# feature functions, will be applied to each g_id
feature_funcs = {'cp': [count],
                    'vol': [np.sum],
                    'lat': np.min,
                    'lon': np.min}
gv = g.partition_nodes('g_id', feature_funcs)
gv.rename(columns={'lat_amin': 'lat',
                         'lon_amin': 'lon'}, inplace=True)
```

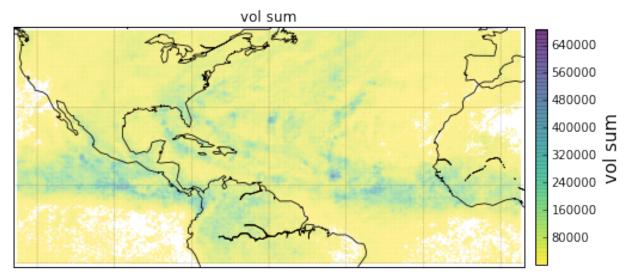
ĺ	n_nodes	cp_count	lat	vol_sum	lon
g_id					
0	2	1	-10.125	10142	-125.125
1	2	1	-9.875	8716	-125.125
2	2	0	-9.625	4372	-125.125
3	2	2	-9.375	5310	-125.125
4	2	2	-9.125	6409	-125.125
115618	2	1	48.875	14319	5.125
115619	1	1	49.125	10129	5.125
115620	2	1	49.375	12826	5.125
115621	2	2	49.625	9117	5.125
115622	2	1	49.875	12101	5.125

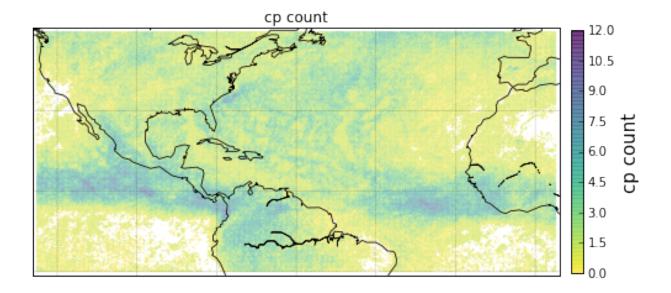
Plot GeoLocational Information

```
cols = { 'n_nodes': gv.n_nodes,
        'vol sum': gv.vol_sum,
        'cp count': gv.cp_count}
for name, col in cols.items():
    # for easy filtering, we create a new DeepGraph instance for
    # each component
   gt = dg.DeepGraph(gv)
    # configure map projection
    kwds_basemap = { 'llcrnrlon': v.lon.min() - 1,
                    'urcrnrlon': v.lon.max() + 1,
                    'llcrnrlat': v.lat.min() - 1,
                    'urcrnrlat': v.lat.max() + 1}
    # configure scatter plots
    kwds_scatter = {'s': 1,
                    'c': col.values,
                    'cmap': 'viridis_r',
                    'alpha': .5,
                    'edgecolors': 'none'}
    # create scatter plot on map
    obj = gt.plot_map(lon='lon', lat='lat',
                      kwds_basemap=kwds_basemap,
                      kwds_scatter=kwds_scatter)
    # configure plots
    obj['m'].drawcoastlines(linewidth=.8)
    obj['m'].drawparallels(range(-50, 50, 20), linewidth=.2)
    obj['m'].drawmeridians(range(0, 360, 20), linewidth=.2)
   obj['ax'].set_title(name)
    # colorbar
```

```
cb = obj['fig'].colorbar(obj['pc'], fraction=.022, pad=.02)
cb.set_label('{}'.format(name), fontsize=15)
```







Geographical Locations and Families

In order to create the intersection partition of geographical locations and families, we first need to append a family membership column to ${\rm v}$

```
# create F col
v['F'] = np.ones(len(v), dtype=int) * -1
gcpv = cpv.groupby('F')
it = gcpv.apply(lambda x: x.index.values)
for F in range(len(it)):
    cp_index = v.cp.isin(it.iloc[F])
    v.loc[cp_index, 'F'] = F
```

Then we create the intersection partition

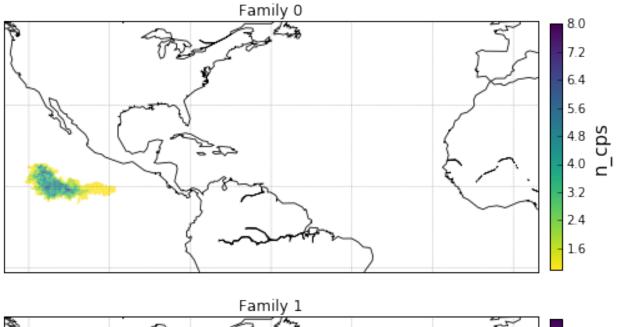
which looks like this

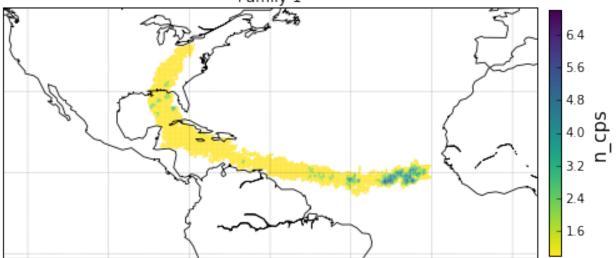
print(fgv)

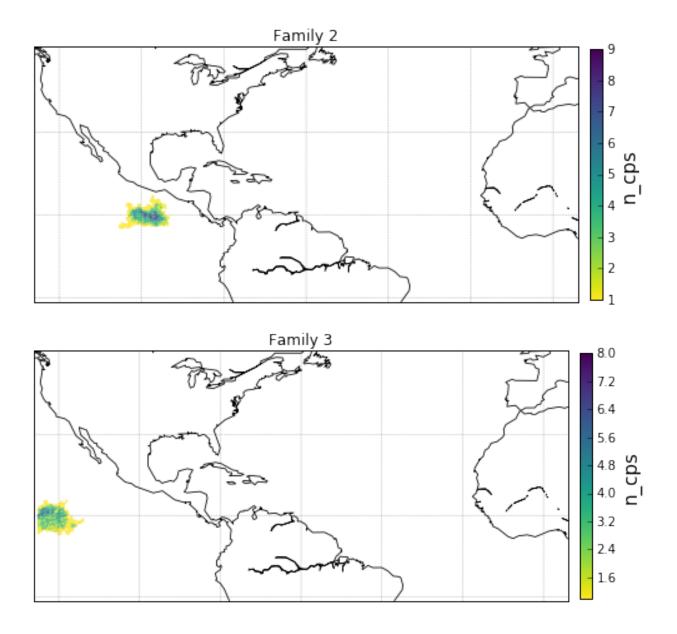
$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$			n_nodes	n_cp_nodes	lat	vol_sum	lon
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	F	g_id					
2 2 1 -9.625 4372 -125.125 3 2 2 -9.375 5310 -125.125 4 2 2 -9.125 6409 -125.125 998 26685 1 1 -8.875 593 -93.625 26686 1 1 -8.625 411 -93.625 26887 1 1 -9.375 364 -93.375 26888 1 1 -9.125 478 -93.375	-1	0	2	2	-10.125	10142	-125.125
$\begin{array}{cccccccccccccccccccccccccccccccccccc$		1	2	2	-9.875	8716	-125.125
4 2 2 -9.125 6409 -125.125 998 26685 1 1 -8.875 593 -93.625 26686 1 1 -8.625 411 -93.625 26887 1 1 -9.375 364 -93.375 26888 1 1 -9.125 478 -93.375		2	2	1	-9.625	4372	-125.125
998 26685 1 1 -8.875 593 -93.625 26686 1 1 -8.625 411 -93.625 26887 1 1 -9.375 364 -93.375 26888 1 1 -9.125 478 -93.375		3	2	2	-9.375	5310	-125.125
998 26685 1 1 -8.875 593 -93.625 26686 1 1 -8.625 411 -93.625 26887 1 1 -9.375 364 -93.375 26888 1 1 -9.125 478 -93.375		4	2	2	-9.125	6409	-125.125
26686 1 1 -8.625 411 -93.625 26887 1 1 -9.375 364 -93.375 26888 1 1 -9.125 478 -93.375							
2688711-9.375364-93.3752688811-9.125478-93.375	998	26685	1	1	-8.875	593	-93.625
26888 1 1 -9.125 478 -93.375		26686	1	1	-8.625	411	-93.625
		26887	1	1	-9.375	364	-93.375
26889 1 1 -8.875 456 -93.375		26888	1	1	-9.125	478	-93.375
		26889	1	1	-8.875	456	-93.375
	[186	903 row	s x 5 col	umns]			

Plot Family Information

```
families = [0, 1, 2, 3]
for F in families:
    # for easy filtering, we create a new DeepGraph instance for
    # each component
   gt = dg.DeepGraph(fgv.loc[F])
    # configure map projection
   kwds_basemap = {'llcrnrlon': v.lon.min() - 1,
                    'urcrnrlon': v.lon.max() + 1,
                    'llcrnrlat': v.lat.min() - 1,
                    'urcrnrlat': v.lat.max() + 1}
    # configure scatter plots
   kwds_scatter = {'s': 1,
                    'c': gt.v.n_cp_nodes.values,
                    'cmap': 'viridis_r',
                    'edgecolors': 'none'}
    # create scatter plot on map
   obj = gt.plot_map(
       lat='lat', lon='lon',
       kwds_basemap=kwds_basemap, kwds_scatter=kwds_scatter)
    # configure plots
   obj['m'].drawcoastlines(linewidth=.8)
   obj['m'].drawparallels(range(-50, 50, 20), linewidth=.2)
   obj['m'].drawmeridians(range(0, 360, 20), linewidth=.2)
   cb = obj['fig'].colorbar(obj['pc'], fraction=.022, pad=.02)
   cb.set_label('n_cps', fontsize=15)
   obj['ax'].set_title('Family {}'.format(F))
```







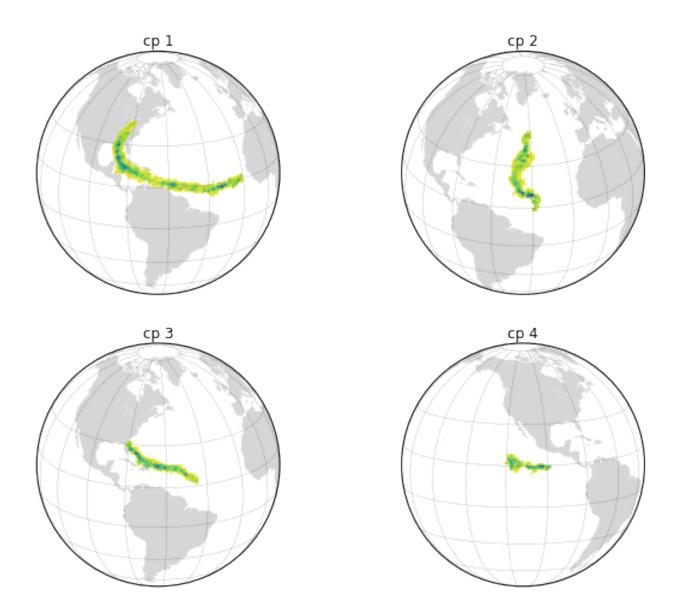
Geographical Locations and Components

print(gcpv)

		n_nodes	lat	vol_sum	lon
ср	g_id				
0	0	1	-10.125	5071	-125.125
	1	1	-9.875	4415	-125.125
	2	2	-9.625	4372	-125.125
	6	3	-8.375	1026	-125.125
	7	1	-8.125	594	-125.125
33167	112117	1	9.375	24618	0.625
33168	100613	1	6.625	11450	-13.625
	100614	1	6.875	12706	-13.625
33169	98523	1	15.375	31057	-16.125
	98524	1	15.625	15741	-16.125
[2873	01 rows	x 4 colum	nns]		

Plot Component Information

```
# select the components to plot
comps = [1, 2, 3, 4]
fig, axs = plt.subplots(2, 2, figsize=[10,8])
axs = axs.flatten()
for comp, ax in zip(comps, axs):
    # for easy filtering, we create a new DeepGraph instance for
    # each component
   gt = dg.DeepGraph(gcpv[gcpv.index.get_level_values('cp') == comp])
    # configure map projection
   kwds_basemap = { 'projection': 'ortho',
                    'lon_0': cpv.loc[comp].lon_mean,
                    'lat_0': cpv.loc[comp].lat_mean,
                    'resolution': 'c'}
    # configure scatter plots
   kwds_scatter = {'s': .5,
                    'c': gt.v.vol_sum.values,
                    'cmap': 'viridis_r',
                    'edgecolors': 'none'}
    # create scatter plot on map
   obj = gt.plot_map(lon='lon', lat='lat',
                      kwds_basemap=kwds_basemap,
                      kwds_scatter=kwds_scatter,
                      ax=ax)
    # configure plots
   obj['m'].fillcontinents(color='0.2', zorder=0, alpha=.2)
   obj['m'].drawparallels(range(-50, 50, 20), linewidth=.2)
   obj['m'].drawmeridians(range(0, 360, 20), linewidth=.2)
   obj['ax'].set_title('cp {}'.format(comp))
```



1.3.4 From Multilayer Networks to Deep Graphs

[ipython notebook] [python script]

In this tutorial we exemplify the representation of multilayer networks (MLNs) by deep graphs and demonstrate some of the advantages of deepgraph's network representation.

We start by converting the Noordin Top Terrorist MLN into a graph g - comprised of two DataFrames, a node table g. v and an edge table g. e - that corresponds to the supra-graph representation of the multilayer network.

We then partition the graph g by the information attributed to its layers, resulting in different supergraphs on the partition lattice of g that correspond to different representations of a MLN (including its tensor representation).

In the next part, we demonstrate how additional information that might be at hand or computed during the analysis can be used to induce further supergraphs, or metaphorically speaking, how additional information corresponds to "hidden layers" of a MLN.

Finally, we briefly show how to use the nodes' properties to partition the edges of a MLN.

** References **

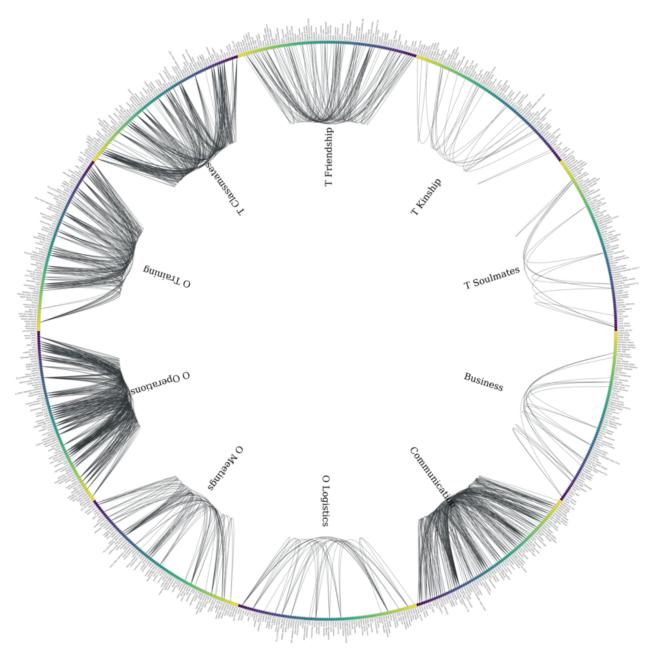
For a short summary of the multilayer network representation, see Appendix C of the Deep Graphs paper.

For a more in-depth introduction to MLNs, I recommend the following papers:

- Multilayer Networks (review paper of MLNs)
- The Structure and Dynamics of Multilayer Networks (review paper of MLNs)
- Mathematical Formulation of Multilayer Networks (tensor formalism for MLNs)

For a discussion of how Deep Graphs relates to the multilayer network representation, see Sec. IV B and Appendix D of the Deep Graphs paper.

The Noordin Top Terrorist Data



[high-res version] [python plot script]

The data we use in this tutorial is the Noordin Top Terrorist Network, which has previously been represented as a multilayer network (e.g., http://arxiv.org/abs/1308.3182)

It includes relational data on 79 Indonesian terrorists belonging to the so-called Noordin Top Terrorist Network.

For information about the individual's attributes and their relations, see http://www.thearda.com/archive/files/ codebooks/origCB/Noordin%20Subset%20Codebook.pdf and http://arxiv.org/pdf/1308.3182v3.pdf.

Preprocessing

We download the data from here, and process it into two pandas DataFrames, a node table and an edge table. The preprocessing is quite lengthy, so you might want to proceed directly to the *next section*.

First of all, we need to import some packages

```
# data i/o
import os
import subprocess
import zipfile
# for plots
import matplotlib.pyplot as plt
# the usual
import numpy as np
import pandas as pd
import deepgraph as dg
# notebook display
%matplotlib inline
pd.options.display.max_rows = 10
pd.set_option('expand_frame_repr', False)
```

Preprocessing the Nodes

```
# there are 10 layers and 79 nodes on each layer
v = pd.concat(10 \star [v])
# add "aspect" as column to v
layer_names = ['Business', 'Communication', 'O Logistics', 'O Meetings',
               'O Operations', 'O Training', 'T Classmates', 'T Friendship',
               'T Kinship', 'T Soulmates']
layers = [[name] *79 for name in layer_names]
layers = [item for sublist in layers for item in sublist]
v['layer'] = layers
# set unique node index
v.reset_index(inplace=True)
v.rename(columns={'index': 'V_N'}, inplace=True)
# swap columns
cols = list(v)
cols[1], cols[10] = cols[10], cols[1]
v = v[cols]
# get rid of the attribute columns for demonstrational purposes,
# will be inserted again later
v, vinfo = v.iloc[:, :2], v.iloc[:, 2:]
```

Preprocessing the Edges

```
# paj file containing edges for different layers
get_paj = ("wget -0 tmp/terrorists.paj "
           "https://sites.google.com/site/sfeverton18/"
           "research/appendix-1/Noordin%20Subset%20%28Pajek%29.paj?"
           "attredirects=0&d=1")
subprocess.call(get_paj.split())
# get data blocks from paj file
with open('tmp/terrorists.paj') as txtfile:
   comments = []
   data = []
    part = []
    for line in txtfile:
        if line.startswith('*'):
            # comment lines
            comment = line
            comments.append(comment)
            if part:
                data.append(part)
                part = []
        else:
            # vertices
            if comment.startswith('*Vertices') and len(line.split()) > 1:
                sublist = line.split('"')
                sublist = sublist[:2] + sublist[-1].split()
                part.append(sublist)
            # edges or partitions
            elif not line.isspace():
```

```
part.append(line.split())
    # append last block
   data.append(part)
# extract edge tables from data blocks
ecomments = []
eparts = []
for i, c in enumerate(comments):
   if c.startswith('*Network'):
        del data[0]
   elif c.startswith('*Partition'):
        del data[0]
    elif c.startswith('*Vector'):
        del data[0]
    elif c.startswith('*Arcs') or c.startswith('*Edges'):
        ecomments.append(c)
        eparts.append(data.pop(0))
# layer data parts (indices found manually via comments)
inds = [11, 10, 5, 6, 7, 8, 0, 1, 2, 3]
eparts = [eparts[ind] for ind in inds]
# convert to DataFrames
layer_frames = []
for name, epart in zip(layer_names, eparts):
   frame = pd.DataFrame(epart, dtype=np.int16)
    # get rid of self-loops, bidirectional edges
   frame = frame[frame[0] < frame[1]]</pre>
    # rename columns
   frame.rename(columns={0: 's', 1: 't', 2: name}, inplace=True)
   frame['s'] -= 1
    frame['t'] -= 1
    layer_frames.append(frame)
# set indices
for i, e in enumerate(layer_frames):
   e['s'] += i*79
   e['t'] += i*79
   e.set_index(['s', 't'], inplace=True)
# concat the layers
e = pd.concat(layer_frames)
# edge table as described in the paper
e_paper = e.copy()
```

```
# alternative representation of e
e['type'] = 0
e['weight'] = 0
for layer in layer_names:
   where = e[layer].notnull()
   e.loc[where, 'type'] = layer
   e.loc[where, 'weight'] = e.loc[where, layer]
e = e[['type', 'weight']]
```

DeepGraph's Supra-Graph Representation of a MLN, G = (V, E)

Above, we have processed the downloaded data into a node table v and an edge table e, that correspond to the supragraph representation of a multilayer network. This is the preferred representation of a MLN by a deep graph, since all other representations are entailed in the supra-graph's partition lattice, as we will demonstrate below.

```
g = dg.DeepGraph(v, e)
print(g)
```

<DeepGraph object, with n=790 node(s) and m=1014 edge(s) at 0x7fb8e13499e8>

Let's have a look at the node table first

prin	c (g.v))	
	V_N		layer
0	0		Business
1	1		Business
2	2		Business
3	3		Business
4	4		Business
785	74	Т	Soulmates
786	75	Т	Soulmates
787	76	Т	Soulmates
788	77	Т	Soulmates
789	78	Т	Soulmates
[790	rows	Х	2 columns]

As you can see, there are 790 nodes in total. Each of the 10 layers,

print(g.v.layer.unique())

```
['Business' 'Communication' 'O Logistics' 'O Meetings' 'O Operations'
'O Training' 'T Classmates' 'T Friendship' 'T Kinship' 'T Soulmates']
```

is comprised of 79 nodes. Every node has a feature of type V_N , indicating the individual the node belongs to, and a feature of type layer, corresponding to the layer the node belongs to. Each of the 790 nodes corresponds to a node-layer of the MLN representation of this data.

The edge table,

print(g.e)

```
type weight
    t
S
9
   67
                         2.0
            Business
    69
                         1.0
            Business
    77
                         1.0
            Business
11
   61
            Business
                         1.0
20
   59
            Business
                         1.0
                 . . .
                         . . .
. . .
733 769 T Soulmates
                         1.0
755 769 T Soulmates
                         1.0
    787 T Soulmates
                         1.0
```

```
771 788 T Soulmates 1.0
783 788 T Soulmates 1.0
[1014 rows x 2 columns]
```

is comprised of 1014 edges between the nodes in v. Each edge has two relations. The first relation (of type type) is determined by the tuple of features $(layer_i, layer_j)$ of the adjacent nodes V_i and V_j . The second relation (of type weight) indicates the "weight" of the connection.

This representation of the edges of a MLN deviates from the one you can find in the paper, which is described in the *last section*.

There are 10 types of relations in the above edge table

```
g.e['type'].unique()
```

which - in the case of this data set - correspond to the layers of the nodes. This is due to the fact that there are no inter-layer connections in the Noordin Top Terrorist Network (such as, e.g., an edge from layer Business to layer Communication would be). The edges here are all (undirected) intra-layer edges (e.g., Business \rightarrow Business, Operations \rightarrow Operations).

To see how the edges are distributed among the different types, you can simply type

```
g.e['type'].value_counts()
```

```
0 Operations
               267
Communication 200
T Classmates
               175
O Training
              147
T Friendship
               91
               63
O Meetings
O Logistics
               29
T Kinship
                16
Business
                15
T Soulmates
                11
Name: type, dtype: int64
```

Let's have a look at how many "actors" (nodes with at least one connection) there are within each layer

```
# append degree
gtg = g.return_gt_graph()
g.v['deg'] = gtg.degree_property_map('total').a
# how many "actors" are there per layer?
g.v[g.v.deg != 0].groupby('layer').size()
```

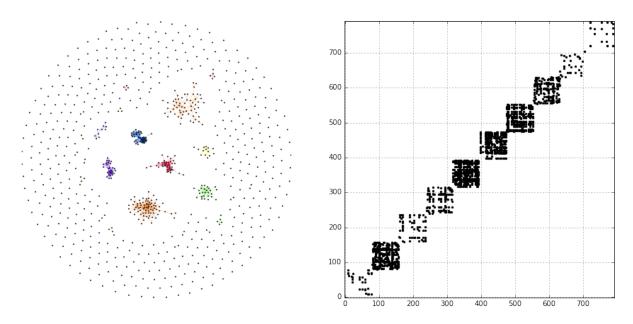
layer Business 13 Communication 74 O Logistics 16 O Meetings 26 O Operations 39

0	Training	38					
Т	Classmates	39					
Т	Friendship	61					
Т	Kinship	24					
Т	Soulmates	9					
dt	dtype: int64						

For the purpose of this tutorial, the fact that the Noordin Top Terrorist Network is a MLN with only one aspect, and without inter-layer edges, is of little importance. The generalization of what we're showing in the following to more general MLNs is straight-forward (and explained in detail in Appendix D of the paper).

Let's illustrate the supra-graph representation of this MLN by a plot

```
# create graph_tool graph for layout
import graph_tool.draw as gtd
gtg = g.return_gt_graph()
gtq.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg, gamma=.5)
pos = pos.get_2d_array([0, 1])
g.v['x'] = pos[0]
g.v['y'] = pos[1]
# configure nodes
kwds_scatter = {'s': 1,
                'c': 'k'}
# configure edges
kwds_guiver = { 'headwidth': 1,
               'alpha': .3,
               'cmap': 'prism'}
# color by type
C = g.e.groupby('type').grouper.group_info[0]
# plot
fig, ax = plt.subplots(1, 2, figsize=(15, 7))
g.plot_2d('x', 'y', edges=True, C=C,
          kwds_scatter=kwds_scatter,
          kwds_quiver=kwds_quiver, ax=ax[0])
# turn axis off, set x/y-lim
ax[0].axis('off')
ax[0].set_xlim((g.v.x.min() - 1, g.v.x.max() + 1))
ax[0].set_ylim((g.v.y.min() - 1, g.v.y.max() + 1))
# plot adjacency matrix
adj = g.return_cs_graph().todense()
adj = adj + adj.T
inds = np.where(adj != 0)
ax[1].scatter(inds[0], inds[1], c='k', marker='.')
ax[1].grid()
ax[1].set_xlim(-1, 791)
ax[1].set_ylim(-1,791)
```



The supra-graph representation of a MLN is by itself a powerful representation and exploitable in various ways (see, e.g., section 2.3 of this paper). However, in the following, we will demonstrate how to use the additional information attributed to the layers of the MLN, in order to "structure" and partition the MLN into different representations.

Redistributing Information on the Partition Lattice of the MLN

Based on the types of features V_N and layer, we can now redistribute the information contained in the supra-graph g. This redistribution allows for several representations of the graph, which we will demonstrate in the following.

The SuperGraph $G^L = (V^L, E^L)$

Partitioning by the type of feature layer leads to the supergraph $G^L = (V^L, E^L)$, where every supernode $V_{iL}^L \in V^L$ corresponds to a distinct layer, encompassing all its respective nodes. Superedges $E_{i^L,j^L}^L \in E^L$ with either $i^L = j^L$ or $i^L \neq j^L$ correspond to collections of intra- and inter-layer edges of the MLN, respectively.

<DeepGraph object, with n=10 node(s) and m=10 edge(s) at 0x7fb8e1349c50>

print(lg.v)

	n_nodes
layer	
Business	79
Communication	79
O Logistics	79
O Meetings	79
O Operations	79

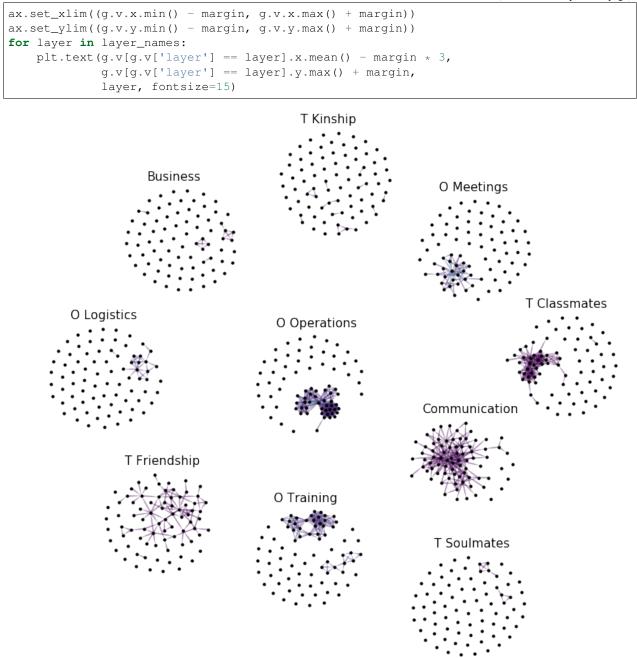
Training	79
Classmates	79
Friendship	79
Kinship	79
Soulmates	79
	Training Classmates Friendship Kinship Soulmates

print(lg.e)

		n_edges	weight_sum	weight_mean	weight_std
layer_s	layer_t				
Business	Business	15	16.0	1.066667	0.258199
Communication	Communication	200	200.0	1.000000	0.00000
O Logistics	O Logistics	29	58.0	2.00000	0.00000
O Meetings	O Meetings	63	170.0	2.698413	1.612801
O Operations	O Operations	267	574.0	2.149813	0.699107
O Training	O Training	147	334.0	2.272109	0.763534
T Classmates	T Classmates	175	175.0	1.000000	0.00000
T Friendship	T Friendship	91	91.0	1.000000	0.00000
T Kinship	T Kinship	16	16.0	1.000000	0.00000
T Soulmates	T Soulmates	11	11.0	1.000000	0.00000

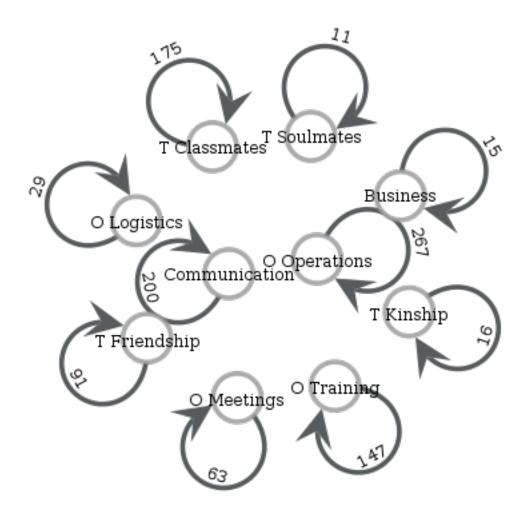
Let's plot the graph g grouped by its layers.

```
# append layer_id to group nodes by layers
g.v['layer_id'] = g.v.groupby('layer').grouper.group_info[0].astype(np.int32)
# create graph_tool graph object
gtg = g.return_gt_graph(features=['layer_id'])
gtg.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg, groups=gtg.vp['layer_id'], mu=.15)
pos = pos.get_2d_array([0, 1])
g.v['x'] = pos[0]
g.v['y'] = pos[1]
# configure nodes
kwds_scatter = {'s': 10,
                'c': 'k'}
# configure edges
kwds_quiver = { 'headwidth': 1,
               'alpha': .4,
               'cmap': 'viridis'}
# color by weight
C = q.e.weight.values
# plot
fig, ax = plt.subplots(figsize=(12, 12))
obj = g.plot_2d('x', 'y', edges=True, C=C,
          kwds_scatter=kwds_scatter,
          kwds_quiver=kwds_quiver, ax=ax)
\ensuremath{\#} turn axis off, set x/y-lim and name layers
ax.axis('off')
margin = 10
```



We can also plot the supergraph $G^L = (V^L, E^L)$

```
edge_text=gtg.ep['n_edges'],
inline=True, fit_view=.8,
output_size=(400,400))
```



The SuperGraph $G^N = (V^N, E^N)$

Partitioning by the type of feature V_N leads to the supergraph $G^N = (V^N, E^N)$, where each supernode $V_{i^N}^N \in V^N$ corresponds to a node of the MLN. Superedges $E_{i^N j^N}^N \in E^N$ with $i^N = j^N$ correspond to the coupling edges of a MLN.

```
# partition by MLN's node indices
nv, ne, gv, ge = g.partition_graph('V_N', return_gve=True)
# for each superedge, get types of edges and their weights
def type_weights(group):
    index = group['type'].values
    data = group['weight'].values
```

```
return pd.Series(data=data, index=index)
ne_weights = ge.apply(type_weights).unstack()
ne = pd.concat((ne, ne_weights), axis=1)
# create graph
ng = dg.DeepGraph(nv, ne)
ng
```

<DeepGraph object, with n=79 node(s) and m=623 edge(s) at 0x7fb8d1da8b70>

print(ng.v)

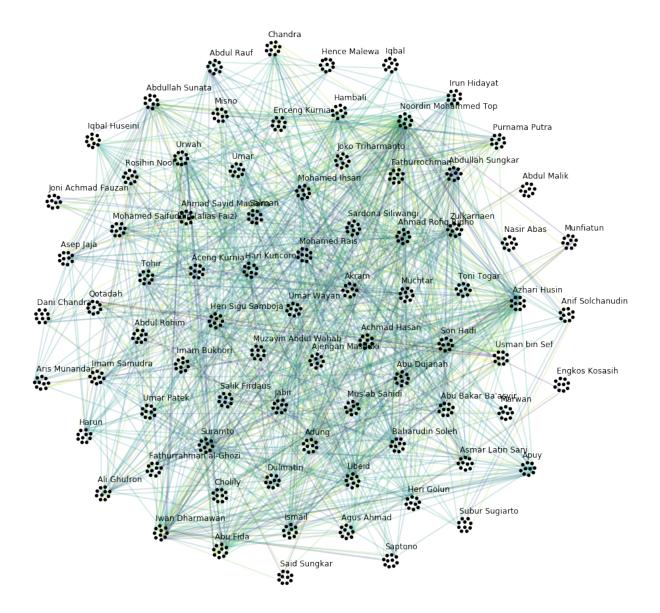
print(ng.e)

n_edgesBusinessCommunicationOLogisticsOMeetingsOOperations.u-OTTClassmatesTFriendshipTKinshipTSoulmatesv_N_sV_N_tVVVVVVVVVV0153NaN1.02.0NaNNaNNaNV141NaNNaNNaNNaNNaNNaNNaNV141NaNNaNNaNNaNNaNNaNNaNV141NaNNaNNaNNaNNaNNaNNaNV141.0NaNNaNNaNNaNNaNNaNNaNV161NaNNaNNaNNaNNaNNaNNaNNaNNaNV211NaNNaNNaNNaNNaNNaNNaNNaNVV14NaNNaNNaNNaNNaNNaNNaNNaNVV15734NaNNaNNaNNaNNaNNaNNaNNaNVV14NaNNaNNaNNaNNaNNaNNaNNaNVVV161.0NaNNaNNaNNaNNaNNaNNaNVVV17734NaNNaNN			n edges	Bugir	1000	Communica	tion	0 1.00	istics	0 Maa	tings	0 Operations
V_{N} s V_{N} t V V_{N} t V_{N} t V V_{N} t V V_{N} t V V V V_{N} t V V V V V V V V V		Tuning	-					-			CINGS	
0153NAN1.02.0NANNANNAN1NANNANNANNANNANNANNANNANNANNAN141NAN<		-	I CLASS	mates	1 1	riendsnip	IVI	nsnip	I SOUL	lliates		
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[623 rows x 11 columns]

```
Let's plot the graph g grouped by V_N.
```

```
# create graph_tool graph object
g.v['V_N'] = g.v['V_N'].astype(np.int32)  # sfpd only takes int32
g_tmp = dg.DeepGraph(v)
gtg = g_tmp.return_gt_graph(features='V_N')
gtg.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg, groups=gtg.vp['V_N'], mu=.3, gamma=.01)
pos = pos.get_2d_array([0, 1])
g.v['x'] = pos[0]
g.v['y'] = pos[1]
# configure nodes
kwds_scatter = { 'c': 'k' }
# configure edges
kwds_quiver = { 'headwidth': 1,
               'alpha': .2,
               'cmap': 'viridis_r'}
# color by type
C = g.e.groupby('type').grouper.group_info[0]
# plot
fig, ax = plt.subplots(figsize=(15,15))
g.plot_2d('x', 'y', edges=True,
          kwds_scatter=kwds_scatter, C=C,
          kwds_quiver=kwds_quiver, ax=ax)
# turn axis off, set x/y-lim and name nodes
name_dic = {i: name for i, name in enumerate(vinfo.iloc[:79].Name)}
ax.axis('off')
ax.set_xlim((g.v.x.min() - 1, g.v.x.max() + 1))
ax.set_ylim((g.v.y.min() - 1, g.v.y.max() + 1))
for node in g.v['V_N'].unique():
   plt.text(g.v[g.v['V_N'] == node].x.mean() - 1,
             g.v[g.v['V_N'] == node].y.max() + 1,
             name_dic[node], fontsize=12)
```

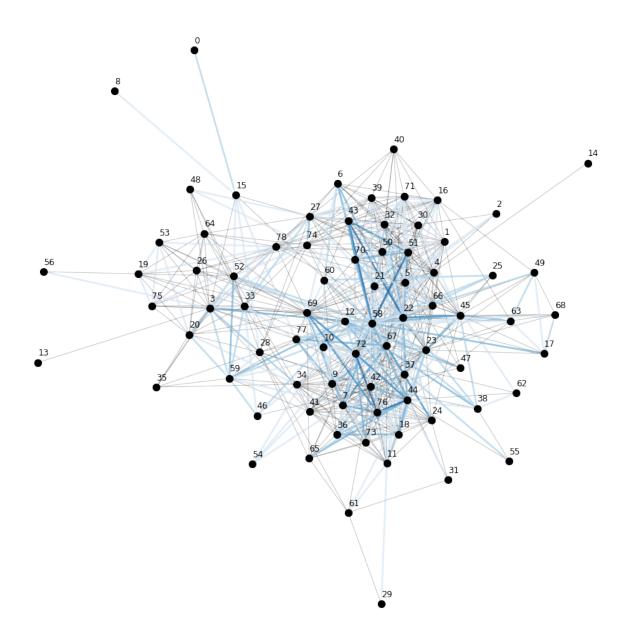


Let's also plot the supergraph $G^N = (V^N, E^N)$, where the color of the superedges corresponds to the number of edges within the respective superedge.

```
# get rid of isolated node for nicer layout
ng.v.drop(57, inplace=True, errors='ignore')
# create graph_tool graph object
gtg = ng.return_gt_graph(features=True, relations='n_edges')
gtg.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg)
pos = pos.get_2d_array([0, 1])
ng.v['x'] = pos[0]
ng.v['y'] = pos[1]
# configure nodes
kwds_scatter = {'s': 100,
```

```
'c': 'k'}
```

```
# configure edges
# split edges with only one type of connection
C_split_0 = ng.e['n_edges'].values.copy()
C_split_0[C_split_0 == 1] = 0
# edges with one type of connection
kwds_quiver_0 = {'alpha': .3,
                 'width': .001}
# edges with more than one type
kwds_quiver = { 'headwidth': 1,
               'width': .003,
               'alpha': .7,
               'cmap': 'Blues',
               'clim': (1, ng.e.n_edges.max())}
# create plot
fig, ax = plt.subplots(figsize=(15, 15))
ng.plot_2d('x', 'y', edges=True, C_split_0=C_split_0,
           kwds_scatter=kwds_scatter, kwds_quiver_0=kwds_quiver_0,
           kwds_quiver=kwds_quiver, ax=ax)
# turn axis off, set x/y-lim and name nodes
ax.axis('off')
ax.set_xlim(ng.v.x.min() - 1, ng.v.x.max() + 1)
ax.set_ylim(ng.v.y.min() - 1, ng.v.y.max() + 1)
for i in ng.v.index:
   plt.text(ng.v.at[i, 'x'], ng.v.at[i, 'y'] + .3, i, fontsize=12)
```



The Tensor-Like Representation $G^{NL} = (V^{NL}, E^{NL})$

Considering only the information attributed to the layers of the MLN, and the fact that this MLN has just one aspect, there is only one more supergraph we can create of g. It is given by creating the intersection partition (see section III E of the Deep Graphs paper) of the types of features V_N and layer. The resulting supergraph $G^{N \cdot L} = (V^{N \cdot L}, E^{N \cdot L})$ corresponds one to one to the graph G = (V, E), and therefore to the supra-graph representation of the MLN. The only difference is the indexing, which is tensor-like for the supergraph $G^{N \cdot L}$.

```
# partition the graph
relation_funcs = {'type': 'sum', 'weight': 'sum'} # just to transfer relations
nlv, nle = g.partition_graph(['V_N', 'layer'], relation_funcs=relation_funcs)
nlg = dg.DeepGraph(nlv, nle)
nlg
```

<DeepGraph object, with n=790 node(s) and m=1014 edge(s) at 0x7fb8d5325550>

print(nlg.v)

```
n_nodes
V_N layer
0 Business
                     1
  Business
Communication
                     1
                     1
   O Logistics
   0 Meetings
                     1
   0 Operations
                    1
• • •
                    . . .
                  _
1
78 O Training
   T Classmates
   T Friendship
                    1
   T Kinship
                     1
                1
   T Soulmates
[790 rows x 1 columns]
```

print(nlg.e)

				n_edges	weight	type
V_N_s	layer_s	V_N_t	layer_t			
0	Communication	15	Communication	1	1.0	Communication
	O Logistics	15	O Logistics	1	2.0	O Logistics
	T Kinship	15	T Kinship	1	1.0	T Kinship
1	O Operations	16	O Operations	1	2.0	O Operations
		22	O Operations	1	2.0	O Operations
72	T Soulmates	77	T Soulmates	1	1.0	T Soulmates
'3	O Operations	76	O Operations	1	2.0	O Operations
	O Training	76	O Training	1	2.0	O Training
75	O Training	78	O Training	1	2.0	O Training
	T Friendship	78	T Friendship	1	1.0	T Friendship
5 1014		2				
[1014	rows x 3 colu	mns]				

This tensor-like index allows you to use the advanced indexing features of pandas.

print(nlg.e.loc[2, 'Communication', :, 'Communication'])

				n_edges	weight	type
V_N_s	layer_s	V_N_t	layer_t			
2	Communication	5	Communication	1	1.0	Communication
		12	Communication	1	1.0	Communication
		30	Communication	1	1.0	Communication
		58	Communication	1	1.0	Communication

In the future, we might implement a method to convert this tensor-representation of a MLN to some sparse-tensor data structure (e.g., https://github.com/mnick/scikit-tensor). Another idea is to create an interface to a suitable multilayer network package that implements the measures and models developed particularly for MLNs.

The "Hidden Layers" of a MLN

Partitioning a multilayer network solely based on the information attributed to its layers only gets us this far. If there is more information available, or computed during the analysis [e.g., by statistical measures, network measures or similarity/distance measures (see g.create_edges)], it can be used to induce further supergraphs and reach other elements of the partition lattice of g.

This is what we'll demonstrate here, based on the additional information available about the individual's attributes:

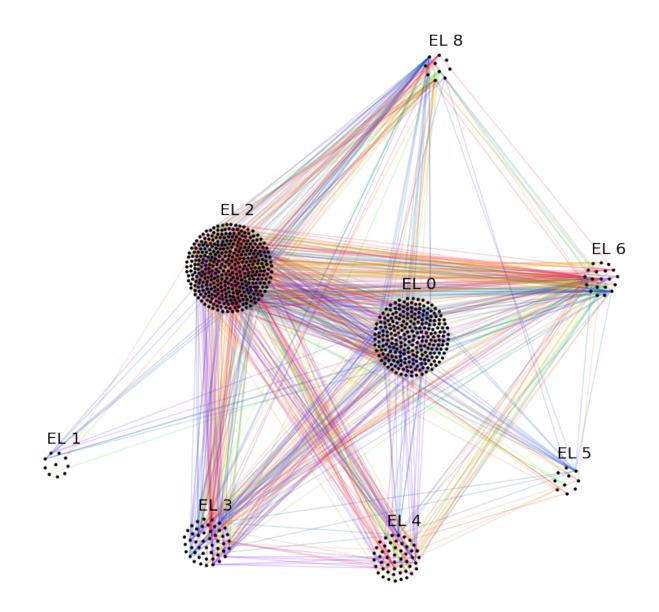
```
print (vinfo)
```

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786	2		1		7	3			
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[790 rows x 9 columns]									

As you can see, there are 9 different attributes associated with each individual, such as their military training, nationality, education level, etc. Let's append this information to the node table, and plot the nodes grouped by their education level.

```
# append node information to g
v = pd.concat((v, vinfo), axis=1)
g = dg.DeepGraph(v, e)
```

```
# create graph_tool graph object
g.v['Education Level'] = g.v['Education Level'].astype(np.int32)
q_tmp = dq.DeepGraph(q.v)
gtg = g_tmp.return_gt_graph(features=['Education Level'])
gtg.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg, groups=gtg.vp['Education Level'], mu=.3, gamma=.1)
pos = pos.get_2d_array([0, 1])
g.v['x'] = pos[0]
g_vv['y'] = pos[1]
# configure nodes
kwds_scatter = {'s': 10,
                'c': 'k'}
# configure edges
kwds_quiver = {'width': 0.002,
               'headwidth': 1,
               'alpha': .2,
               'cmap': 'prism'}
# color by type
C = g.e.groupby('type').grouper.group_info[0]
# plot
fig, ax = plt.subplots(figsize=(13, 12))
obj = g.plot_2d('x', 'y', edges=True,
          kwds_scatter=kwds_scatter, C=C,
          kwds_quiver=kwds_quiver, ax=ax)
# turn axis off, set x/y-lim and name layers
ax.axis('off')
ax.set_xlim((g.v.x.min() - 1, g.v.x.max() + 1))
ax.set_ylim((g.v.y.min() - 1, g.v.y.max() + 1))
for el in g.v['Education Level'].unique():
   plt.text(g.v[g.v['Education Level'] == el].x.mean() - 1,
             g.v[g.v['Education Level'] == el].y.max() + 1,
             'EL {}'.format(el), fontsize=20)
```

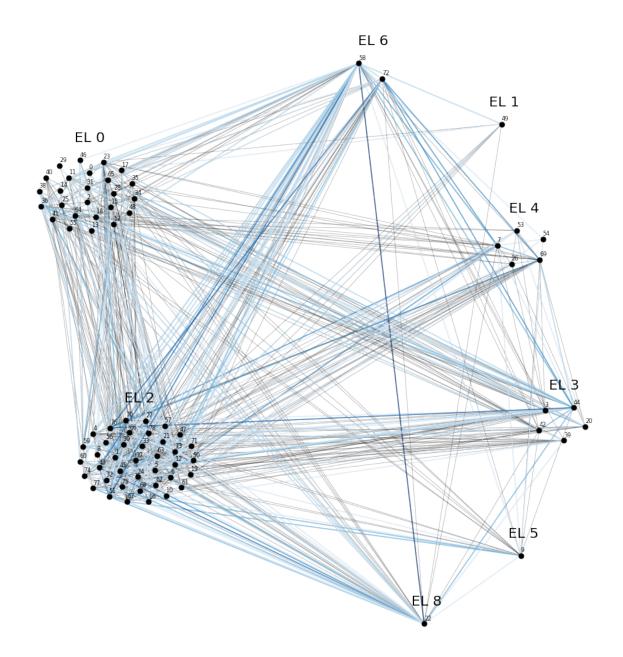


Let's also append the information to the supergraph G^N , and plot this supergraph grouped by education level.

```
# append info to ng.v
ng.v = pd.concat((ng.v, vinfo[:79]), axis=1)
```

```
# create graph_tool graph object
ng.v['Education Level'] = ng.v['Education Level'].astype(np.int32)
g_tmp = dg.DeepGraph(ng.v)
gtg = g_tmp.return_gt_graph(features=['Education Level'])
gtg.set_directed(False)
# get sfdp layout postitions
pos = gtd.sfdp_layout(gtg, groups=gtg.vp['Education Level'], mu=.3, gamma=.01)
pos = pos.get_2d_array([0, 1])
ng.v['x'] = pos[0]
ng.v['y'] = pos[1]
```

```
# configure nodes
kwds_scatter = {'s': 50,
                'c': 'k'}
# configure edges
# split edges with only one type of connection
C_split_0 = ng.e['n_edges'].values.copy()
C_split_0[C_split_0 == 1] = 0
# edges with one type of connection
kwds_quiver_0 = {'alpha': .3,
                 'width': .001}
# edges with more than one type
kwds_quiver = { 'headwidth': 1,
               'width': .002,
               'alpha': .7,
               'cmap': 'Blues',
               'clim': (1, ng.e.n_edges.max()) }
# create plot
fig, ax = plt.subplots(figsize=(15,15))
obj = ng.plot_2d('x', 'y', edges=True, C_split_0=C_split_0,
                 kwds_scatter=kwds_scatter, kwds_quiver_0=kwds_quiver_0,
                 kwds_quiver=kwds_quiver, ax=ax)
# turn axis off, set x/y-lim and name nodes
ax.axis('off')
ax.set_xlim(ng.v.x.min() - 1, ng.v.x.max() + 1)
ax.set_ylim(ng.v.y.min() - 1, ng.v.y.max() + 1)
for i in ng.v.index:
   plt.text(ng.v.at[i, 'x'],
             ng.v.at[i, 'y'] + .2,
             i, fontsize=8)
for el in ng.v['Education Level'].unique():
   plt.text(ng.v[ng.v['Education Level'] == el].x.mean() - .5,
             ng.v[ng.v['Education Level'] == el].y.max() + 1,
             'EL {}'.format(el), fontsize=20)
```



We can now further partition the supergraph G^N into groups with the same education level.

```
df2.index.names = df2.index.names[::-1]
ELne = df1.add(df2, fill_value=0)
# set dtypes
for col in ELne.columns:
    ELne[col] = ELne[col].astype(int)
# find the type of connection most dominant between supernodes
ELne['dominant_type'] = ELne[layer_names].idxmax(axis=1)
# change column order
ELne = ELne[['n_edges'] + ['dominant_type'] + layer_names]
# create graph
ELng = dg.DeepGraph(ELnv, ELne)
ELng
```

<DeepGraph object, with n=8 node(s) and m=30 edge(s) at 0x7fb8d1d245c0>

print(ELng.v)

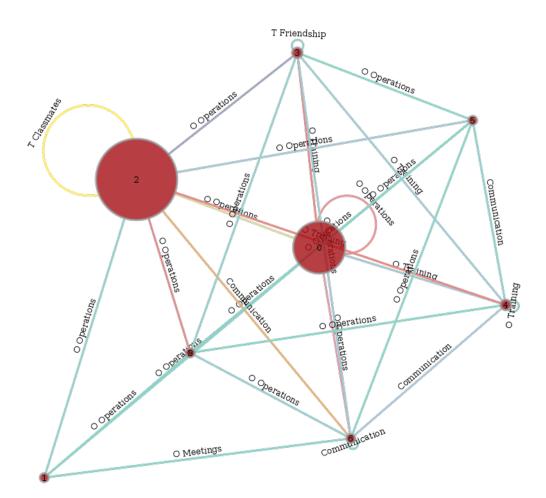
	n_nodes
Education	
0	25
1	1
2	39
3	5
4	5
5	1
6	2
8	1

print(ELng.e)

					n_e	dges	domina	nt_type	Busin	ess	Communica	tion	
⇔ 0	Logi	stics O	Meeti	ngs	0 Operations								-
	-	o T Sou		-	1		2				1		
		-			n Level_t								
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Let's plot the supergraph of education levels, where the node size relates to the number of individuals, edge colors correspond to the number of edges, and edge labels correspond to the most dominant type of connection between nodes.



Partitioning Edges Based on Node Properties

Here, we demonstrate very briefly how to use the additional information of the nodes to perform queries on the edges.

```
# create "undirected" edge table (swap-copy all edges)
g.e = pd.concat((e, e.swaplevel(0,1)))
g.e.sort_index(inplace=True)
```

```
print(g.partition_edges(source_features=['Nationality']))
```

```
n_edges
Nationality_s
```

3 1655
4 351
5 22

			n_edges	
Nationality_s	Military	Training_t		
3	0		185	
	1		51	
	3		847	
	4		60	
	5		115	
• • •				
5	4		3	
	5		1	
	7		1	
	9		1	
	10		1	
100				
[26 rows x 1 0	columns]			

			n_edges	
type	Nationality_s	Military Training_t		
Business	3	0	3	
		3	16	
		4	1	
		9	2	
		10	2	
T Soulmates	3	9	1	
		10	2	
	4	3	3	
		9	1	
		10	3	
[138 rows x	1 columns]			

Alternative Representation of the MLN Edges

The edges of the supra-graph representation as presented in the paper look like this

print(e_paper)

лт				O Logistics Kinship T Sor	-	0 Operations	O Training 🛓
⇒ı s	t	mates i rii			uimates		
9	67	2.0	NaN	NaN	NaN	NaN	NaN 🔒
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\hookrightarrow		NaN	NaN		NaN		NaN				
20	59	1.		NaN		NaN		NaN	NaN	NaN	ш
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As you can see, the edge table is also comprised of 1014 edges between the nodes in v. However, every type of connection get's its own column, where a "nan" value means that an edge does not have a relation of the corresponding type.

What Next

Now that you have an idea of what the DeepGraph package provides, you should investigate the parts of the package most useful for you. See *API Reference* for details.

1.4 API Reference

The API reference summarizes DeepGraph's core class, its methods and the functions subpackage.

1.4.1 The DeepGraph class

<pre>DeepGraph([v, e, supernode_labels_by,])</pre>	The core class of DeepGraph (dg).	
	The core class of 2 cop or up (ug).	

deepgraph.deepgraph.DeepGraph

class DeepGraph (*v*=*None*, *e*=*None*, *supernode_labels_by*=*None*, *superedge_labels_by*=*None*) The core class of DeepGraph (dg).

This class encapsulates the graph representation as pandas.DataFrame objects in its attributes v and e. It can be initialized with a node table v, whose rows represent the nodes of the graph, as well as an edge table e, whose rows represent edges between the nodes.

Given a node table v, it provides methods to iteratively compute pairwise relations between the nodes using arbitrary, user-defined functions. These methods provide arguments to parallelize the computation and control memory consumption (see create_edges and create_edges_ft).

Also provides methods to partition nodes, edges or an entire graph by the graph's properties and labels, and to create common network representations and graph objects of popular Python network packages.

Furthermore, it provides methods to visualize graphs and their properties and to benchmark the graph construction parameters.

Optionally, the convenience parameter $supernode_labels_by$ can be passed, creating supernode labels by enumerating all distinct (tuples of) values of a (multiple) column(s) of v. Superedge labels can be created analogously, by passing the parameter $superedge_labels_by$.

Parameters

- v (pandas.DataFrame or pandas.HDFStore, optional (default=None)) - The node table, a table representation of the nodes of a graph. The index of v must be unique and represents the node indices. The column names of v represent the types of features of the nodes, and each cell represents a feature of a node. Only a reference to the input DataFrame is created, not a copy. May also be a pandas.HDFStore, but only create_edges and create_edges_ft may then be used (so far).
- e (pandas.DataFrame, optional (default=None)) The edge table, a table representation of the edges between the nodes given by v. Its index has to be a pandas. core.index.MultiIndex, whose first level contains the indices of the source nodes, and the second level contains the indices of the target nodes. Each row of e represents an edge, column names of e represent the types of relations of the edges, and each cell in e represents a relation of an edge. Only a reference to the input DataFrame is created, not a copy.
- supernode_labels_by (dict, optional (default=None)) A dictionary whose keys are strings and their values are (lists of) column names of v. Appends a column to v for each key, whose values correspond to supernode labels, enumerating all distinct (tuples of) values of the column(s) given by the dict's value.
- **superedge_labels_by** (*dict*, *optional* (*default=None*)) A dictionary whose keys are strings and their values are (lists of) column names of e. Appends a column to e for each key, whose values correspond to superedge labels enumerating all distinct (tuples of) values of the column(s) given by the dict's value.

v

See Parameters.

Type pandas.DataFrame

е

See Parameters.

Type pandas.DataFrame

n

Property: Number of nodes.

Type int

m

Property: Number of edges.

Type int

f

Property: types of features and number of features of corresponding type.

Type pd.DataFrame

r

Property: types of relations and number of relations of corresponding type.

Type pd.DataFrame

____init___ (v=None, e=None, supernode_labels_by=None, superedge_labels_by=None) Initialize self. See help(type(self)) for accurate signature.

Methods

init([v, e, supernode_labels_by,])	Initialize self.
<pre>append_binning_labels_v(col, col_name[,</pre>	Append a column with binning labels of the values
])	inv[col].
<pre>append_cp([directed, connection, col_name,])</pre>	Append a component membership column to v.
append_datetime_categories_v([col,	Append datetime categories to v.
])	
<pre>create_edges([connectors, selectors,])</pre>	Create an edge table e linking the nodes in v .
<pre>create_edges_ft(ft_feature[, connectors,])</pre>	Create (ft) an edge table e linking the nodes in v .
filter_by_interval_e(col, interval[, end-	Keep only edges in e with relations of type col in
point])	interval.
<pre>filter_by_interval_v(col, interval[, end-</pre>	Keep only nodes in v with features of type col in
point])	interval.
filter_by_values_e(col, values)	Keep only edges in e with relations of type col in
	values.
filter_by_values_v(col,values)	Keep only nodes in v with features of type col in
	values.
<pre>partition_edges([relations,])</pre>	Return a superedge DataFrame se.
<pre>partition_graph(features[, feature_funcs,</pre>	Return supergraph DataFrames sv and se.
])	
<pre>partition_nodes(features[, feature_funcs,</pre>	Return a supernode DataFrame sv.
])	
<pre>plot_2d(x, y[, edges, C, C_split_0,])</pre>	Plot nodes and corresponding edges in 2 dimensions.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,])</pre>	Plot nodes and corresponding edges by groups.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,])</pre>	Plot nodes and corresponding edges by groups. Work in progress!
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile)</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.
plot_2d(x, y[, edges, C, C_split_0,])plot_2d_generator(x, y, by[, edges, C,])plot_3d(x, y, z[, edges, kwds_scatter,])plot_hist(x[, bins, log_bins, density,])plot_logfile(logfile)plot_map(lon, lat[, edges, C, C_split_0,])	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,]) plot_rects_numeric_numeric(yb, yt, xl, xr[,])</pre>	Plot nodes and corresponding edges by groups. Work in progress! Plot a histogram (or pdf) of x. Plot a logfile. Plot nodes and corresponding edges on a basemap. Plot nodes and corresponding edges by groups, on basemaps. Work in progress! Work in progress! Work in progress!
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,]) plot_rects_numeric_numeric(yb, yt, xl,])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix represen-
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,]) plot_rects_numeric_numeric(yb, yt, xl, xr[,]) return_cs_graph([relations, dropna])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix representation(s).
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_rester(label[, time, ax]) plot_rests_label_numeric(label, xl, xr[,]) plot_rests_numeric_numeric(yb, yt, xl, xr[,]) return_cs_graph([relations, dropna]) </pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix represen- tation(s).Return a graph_tool.Graph representation.
plot_2d(x, y[, edges, C, C_split_0,])plot_2d_generator(x, y, by[, edges, C,])plot_3d(x, y, z[, edges, kwds_scatter,])plot_hist(x[, bins, log_bins, density,])plot_logfile(logfile)plot_map(lon, lat[, edges, C, C_split_0,])plot_map_generator(lon, lat, by[, edges, C,])plot_rester(label[, time, ax])plot_rects_label_numeric(label, xl, xr[,])plot_rects_numeric_numeric(yb, yt, xl, xr[,])return_cs_graph([relations, dropna])return_nx_graph([features, relations,])	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix representation(s).Return a graph_tool.Graph representation.Return a networkx.DiGraph representation.
<pre>plot_2d(x, y[, edges, C, C_split_0,]) plot_2d_generator(x, y, by[, edges, C,]) plot_3d(x, y, z[, edges, kwds_scatter,]) plot_hist(x[, bins, log_bins, density,]) plot_logfile(logfile) plot_map(lon, lat[, edges, C, C_split_0,]) plot_map_generator(lon, lat, by[, edges, C,]) plot_raster(label[, time, ax]) plot_rects_label_numeric(label, xl, xr[,]) plot_rects_numeric_numeric(yb, yt, xl, xr[,]) return_cs_graph([features, relations,]) return_nx_graph([features, relations, dropna]) return_nx_multigraph([features, relations, dropna])</pre>	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix represen- tation(s).Return a graph_tool.Graph representation.
plot_2d(x, y[, edges, C, C_split_0,])plot_2d_generator(x, y, by[, edges, C,])plot_3d(x, y, z[, edges, kwds_scatter,])plot_hist(x[, bins, log_bins, density,])plot_logfile(logfile)plot_map(lon, lat[, edges, C, C_split_0,])plot_map_generator(lon, lat, by[, edges, C,])plot_rester(label[, time, ax])plot_rects_label_numeric(label, xl, xr[,])plot_rects_numeric_numeric(yb, yt, xl, xr[,])return_cs_graph([relations, dropna])return_nx_graph([features, relations,])	Plot nodes and corresponding edges by groups.Work in progress!Plot a histogram (or pdf) of x.Plot a logfile.Plot nodes and corresponding edges on a basemap.Plot nodes and corresponding edges by groups, on basemaps.Work in progress!Work in progress!Work in progress!Work in progress!Return scipy.sparse.coo_matrix representation(s).Return a graph_tool.Graph representation.Return a networkx.DiGraph representation.

Attributes

f	Types of features and number of features of corre-
	sponding type.
m	The number of edges
п	The number of nodes
r	Types of relations and number of relations of corre-
	sponding type.

Creating Edges

<pre>DeepGraph.create_edges([connectors,])</pre>	Create an edge table e linking the nodes in v .
<pre>DeepGraph.create_edges_ft(ft_feature[,])</pre>	Create (ft) an edge table e linking the nodes in v .

deepgraph.deepgraph.DeepGraph.create_edges

DeepGraph.**create_edges** (connectors=None, selectors=None, transfer_features=None, r_dtype_dic=None, no_transfer_rs=None, step_size=10000000, from_pos=0, to_pos=None, hdf_key=None, verbose=False, logfile=None) Create an edge table e linking the nodes in v.

This method enables an iterative computation of pairwise relations (edges) between the nodes represented by v. It does so in a flexible, efficient and vectorized fashion, easily parallelizable and with full control over RAM usage.

1. Connectors

The simplest use-case is to define a single connector function acting on a single column of the node table v. For instance, given a node table v

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'time': [0.,2.,9.], 'x': [3.,1.,12.]})
>>> g = dg.DeepGraph(v)
```

>>> g.v
time x
0 0 3
1 2 1
2 9 12

one may define a function

```
>>> def time_difference(time_s, time_t):
... dt = time_t - time_s
... return dt
```

and pass it to create_edges, in order to compute the time difference of each pair of nodes

```
>>> g.create_edges(connectors=time_difference)
```

```
>>> g.e
dt
s t
```

(continues on next page)

(continued from previous page)

0 1	2	2																							
2	9)																							
1 2	2 9 7	7																							

As one can see, the connector function takes column names of v with additional '_s' and '_t' endings (indicating source node values and target node values, respectively) as input, and returns a variable with the computed values. The resulting edge table g.e is indexed by the node indices ('s' and 't', representing source and target node indices, respectively), and has one column ('dt', the name of the returned variable) with the computed values of the given connector. Note that only the upper triangle adjacency matrix is computed, which is always the case. See Notes for further information.

One may also pass a list of functions to connectors, which are then computed in the list's order. Generally, a connector function can take multiple column names of v (with '_s' and/or '_t' appended) as input, as well as already computed relations of former connectors. Also, any connector function may have multiple output variables. Every output variable has to be a 1-dimensional np.ndarray (with arbitrary dtype, including object). The return statement may not contain any operators, only references to each computed relation.

For instance, considering the above example, one may define an additional connector

```
>>> def velocity(dt, x_s, x_t):
... dx = x_t - x_s
... v = dx / dt
... return v, dx
```

and then apply both connectors on v, resulting in

```
>>> g.create_edges(connectors=[time_difference, velocity])
```

```
>>> g.e
    dt dx v
s t
0 1 2 -2 -1.000000
2 9 9 1.000000
1 2 7 11 1.571429
```

2. Selectors

However, one is often only interested in a subset of all possible edges. In order to select edges during the iteration process - based on some conditions on the node's features and their computed relations - one may pass a (list of) selector function(s) to create_edges. For instance, given the above example, one may define a selector

```
>>> def dt_thresh(dt, sources, targets):
... sources = sources[dt > 5]
... targets = targets[dt > 5]
... return sources, targets
```

and apply it in conjunction with the time_difference connector

```
>>> g.create_edges(connectors=time_difference, selectors=dt_thresh)
```

```
>>> g.e
dt
s t
```

(continues on next page)

		(continued from previous page)
0 2	9	
1 2	7	

leaving only edges with a time difference larger than 5.

Every selector function must have sources and targets as input arguments as well as in the return statement. Most generally, they may depend on column names of v (with '_s' and/or '_t' appended) and/or computed relations of connector functions, and/or computed relations of former selector functions. Apart from sources and targets, they may additionally return computed relations. Given this input/output flexibility of selectors, one could in fact compute all required relations, and select any desired subset of edges, with a single selector function. The purpose of splitting connectors and/or selectors, however, is to control the iteration's performance by consecutively computing relations and selecting edges: **hierarchical selection**.

3. Hierarchical Selection

As the algorithm iterates through the chunks of all possible source and target node indices $([0, g.n^*(g.n-1)/2])$, it goes through the list of selectors at each step. If a selector has a relation as input, it must have either been computed by a former selector, or the selector requests its computation by the corresponding connector function in connectors (this connector may not depend on any other not yet computed relations). Once the input relations are computed (if requested), the selector is applied and returns updated indices, which are then passed to the next selector. Hence, with each selector, the indices are reduced and consecutive computation of relations only consider the remaining indices. After all selectors have been applied, the connector functions that have not been requested by any selector are computed (on the final, reduced chunk of node and target indices).

4. Transferring Features

The argument transfer_features, which takes a (list of) column name(s) of v, makes it possible to transfer features of v to the created edge table e

```
>>> g.create_edges(connectors=time_difference,
                     transfer_features=['x', 'time'])
. . .
>>> q.e
     dt.
          time_s time_t
                            X_S
                                 x t
s t
0
 1
      2
               0
                        2
                              3
                                    1
  2
      9
               0
                        9
                              3
                                   12
      7
                        9
1 2
               2
                              1
                                   12
```

If computation time and memory consumption are of no concern, one might skip the remaing paragraphs.

5. Logging

Clearly, the order of the hierarchical selection as described in 3. influences the computation's efficiency. The complexity of a relation's computation and the (expected average) number of deleted edges of a selector should be considered primarily. In order to track and benchmark the iteration process, the progress and time measurements are printed for each iteration step, if verbose is set to True. Furthermore, one may create a logfile (which can also be plot by dg.DeepGraph.plot_logfile) by setting the argument logfile to a string, indicating the file name of the created logfile.

6. Parallelization and Memory Control

The arguments from_pos, to_pos and step_size control the range of processed pairs of nodes and the number of pairs of nodes to process at each iteration step. They may be used for parallel computation and to control RAM usage. See Parameters for details.

It is also possible to initiate dg.DeepGraph with a pandas.HDFStore containing the DataFrame representing the node table. Only the data requested by transfer_features and the user- defined

connectors and selectors at each iteration step is then pulled from the store, which is particularly useful for large node tables and parallel computation. The only requirement is that the node table contained in the store is in table(t) format, not fixed(f) format. For instance, considering the above created node table, one may store it in a hdf file

```
>>> vstore = pd.HDFStore('vstore.h5')
>>> vstore.put('node_table', v, format='t', index=False)
```

initiate a DeepGraph instance with the store

```
>>> g = dg.DeepGraph(vstore)
```

```
>>> g.v
<class 'pandas.io.pytables.HDFStore'>
File path: vstore.h5
/node_table frame_table (typ->appendable,nrows->3,ncols->2,
indexers->[index])
```

and then create edges the same way as if g.v were a DataFrame

```
>>> g.create_edges(connectors=time_difference)
```

```
>>> g.e
dt
s t
0 1 2
2 9
1 2 7
```

In case the store has multiple nodes, hdf_key has to be set to the node corresponding to the node table of the graph.

Also, one may pass a (list of) name(s) of computed relations, no_transfer_rs, which should not be transferred to the created edge table e. This can be advantageous, for instance, if a selector depends on computed relations that are of no further interest.

Furthermore, it is possible to force the dtype of computed relations with the argument r_dtype_dic . The dtype of a relation is then set at each iteration step, but **after** all selectors and connectors were processed.

7. Creating Edges on a Fast Track

If the selection of edges includes a simple distance threshold, i.e. a selector function defined as follows:

```
>>> def ft_selector(x_s, x_t, threshold, sources, targets):
... dx = x_t - x_s
... sources = sources[dx <= threshold]
... targets = targets[dx <= threshold]
... return sources, targets, dx
```

the method create_edges_ft should be considered, since it provides a much faster iteration algorithm.

Parameters

• **connectors** (function or array_like, optional (default=None)) – User defined connector function(s) that compute pairwise relations between the nodes in v. A connector accepts multiple column names of v (with '_s' and/or '_t' appended, indicating source node values and target node values, respectively) as input, as well as already computed relations of former connectors. A connector function may have multiple output variables. Every output variable has to be a 1-dimensional np.ndarray (with arbitrary dtype, including object). See above and dg.functions for examplary connector functions.

- **selectors** (function or array_like, optional (default=None)) User defined selector function(s) that select edges during the iteration process, based on some conditions on the node's features and their computed relations. Every selector function must have sources and targets as input arguments as well as in the return statement. A selector may depend on column names of v (with '_s' and/or '_t' appended) and/or computed relations of connector functions, and/or computed relations of former selector functions. Apart from sources and targets, they may also return computed relations (see connectors). See above, and dg.functions for exemplary selector functions.
- transfer_features (str, int or array_like, optional (default=None)) A (list of) column name(s) of v, indicating which features of v to transfer to e (appending '_s' and '_t' to the column names of e, indicating source and target node features, respectively).
- **r_dtype_dic** (*dict*, *optional* (*default=None*)) A dictionary with names of computed relations of connectors and/or selectors as keys and dtypes as values. Forces the data types of the computed relations in e during the iteration (but **after** all selectors and connectors were processed), otherwise infers them.
- no_transfer_rs (str or array_like, optional (default=None)) Name(s) of computed relations that are not to be transferred to the created edge table e. Can be used to save memory, e.g., if a selector depends on computed relations that are of no interest otherwise.
- **step_size** (*int*, *optional* (*default=1e6*)) The number of pairs of nodes to process at each iteration step. Must be in [1, g.n*(g.n-1)/2]. Its value determines computation speed and memory consumption.
- **from_pos** (*int*, *optional* (*default=0*)) Determines from which pair of nodes to start the iteration process. Must be in [0, g.n*(g.n-1)/2 [. May be used in conjuction with to_pos for parallel computation.
- to_pos (positive integer, optional (default=None)) Determines at which pair of nodes to stop the iteration process (the endpoint is excluded). Must be in [1, g.n*(g.n-1)/2] and larger than from_pos. Defaults to None, which translates to the last pair of nodes, g.n*(g.n-1)/2. May be used in conjunction with from_pos for parallel computation.
- hdf_key (*str*, *optional* (*default=None*)) If you initialized dg. DeepGraph with a pandas.HDFStore and the store has multiple nodes, you must pass the key to the node in the store that corresponds to the node table.
- **verbose** (*bool*, *optional* (*default=False*)) Whether to print information at each step of the iteration process.
- **logfile** (*str*, *optional* (*default=None*)) Create a log-file named by logfile. Contains the time and date of the method's call, the input arguments and time mesaurements for each iteration step. A plot of logfile can be created by dg. DeepGraph.plot_logfile.

Returns e – Set the created edge table e as attribute of dg.DeepGraph.

Return type pd.DataFrame

See also:

```
create_edges_ft()
```

Notes

1. Input and output data types

Since connectors (and selectors) take columns of a pandas DataFrame as input, there are no restrictions on the data types of which pairwise relations are computed. In the most general case, a DataFrame's column has object as dtype, and its values may then be arbitrary Python objects. The same goes for the output variables of connectors (and selectors). The only requirement is that each ouput variable is 1-dimensional.

However, it is also possible to use the values of a column of v as references to arbitrary objects, which may sometimes be more convenient. In case a connector (or selector) needs the node's original indices as input, one may simply copy them to a column, e.g.

>>> v['indices'] = v.index

and then define the connector's (or selector's) input arguments accordingly.

2. Connectors and selectors

The only requirement on connectors and selectors is that their input arguments and return statements are consistent with the column names of v and the passing of computed relations (see above, 3. Hierarchical Selection).

Whatever happens inside the functions is entirely up to the user. This means, for instance, that one may wrap arbitrary functions within a connector (selector), such as optimized C functions or existing functions whose input/output is not consistent with the create_edges method (see, e.g., the methods provided in dg.functions, scipy or scikit learn's sklearn.metrics and sklearn.neighbors. DistanceMetric). One could also store a connector's (selector's) computations directly within the function, or let the function print out any desired information during iteration.

3. Why not compute the full adjacency matrix?

This is due to efficiency. For any asymmetric function (i.e., f(s, t) = f(t, s)), one can always create an additional connector (or output variable) that computes the mirrored values of that function.

deepgraph.deepgraph.DeepGraph.create_edges_ft

DeepGraph.create_edges_ft (ft_feature, connectors=None, selectors=None, transfer_features=None, r_dtype_dic=None, no_transfer_rs=None, min_chunk_size=1000, max_pairs=10000000, from_pos=0, to_pos=None, hdf_key=None, verbose=False, logfile=None)

Create (ft) an edge table e linking the nodes in v.

This method implements the same functionalities as create_edges, with the difference of providing a much quicker iteration algorithm based on a so-called fast-track feature. It is advised to read the docstring of create_edges before this one, since only the differences are explained in the following.

Apart from the hierarchical selection through connectors and selectors as described in the method create_edges (see 1.-3.), this method necessarily includes the (internal) selector function

```
>>> def ft_selector(ftf_s, ftf_t, ftt, sources, targets):
... ft_r = ftf_t - ftf_s
... sources = sources[ft_r <= ftt]
... targets = targets[ft_r <= ftt]
... return sources, targets, ft_r</pre>
```

where ftf is the fast-track feature (a column name of v), ftt the fast-track threshold (a positive number), and ft_r the computed fast-track relation. The argument ft_feature, which has to be a tuple (ftf, ftt), determines these variables.

1. The Fast-Track Feature

The simplest use-case, therefore, is to only pass ft_feature. For instance, given a node table

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'time': [-3.6,-1.1,1.4,4., 6.3],
... 'x': [-3.,3.,1.,12.,7.]})
>>> g = dg.DeepGraph(v)
```

>>> q.v time Х 0 -3.6 -3 -1.1 3 1 2 1.4 1 3 4.0 12 4 6.3 7

one may create and select edges by

```
>>> g.create_edges_ft(ft_feature=('time', 5))
>>> g.e
     ft r
s t
      2.5
0 1
      5.0
 2
1 2
      2.5
2
 3
      2.6
  4
      4.9
34
      2.3
```

leaving only edges with a time difference smaller than (or equal to) ftt = 5. Note that the node table always has to be sorted by the fast-track feature. This is due to the fact that the algorithm only processes pairs of nodes whose fast-track relation is smaller than (or equal to) the fast-track threshold, and the (pre)determination of these pairs relies on a sorted DataFrame.

2. Hierarchical Selection

Additionally, one may define connectors and selectors as described in create_edges (see 1.-3.). Per default, the (internal) fast-track selector is applied first. It's order of application, however, may be determined by inserting the string 'ft_selector' in the desired position of the list of selectors.

The remaining arguments are as described in create_edges, apart from min_chunk_size, max_pairs, from_pos and to_pos. If computation time and/or memory consumption are a concern, one may therefore read the remaining paragraph.

3. Parallelization and Memory Control on a FastTrack

At each iteration step, the algorithm takes a number of nodes ($n = min_chunk_size$, per default n=1000) and computes the fast track relation (distance) between the last node and the first node, $d_ftf = ftf_last - ftf_first$. In case $d_ftf > ftt$, all nodes with a fast- track feature < ftf_last - ftt are considered source nodes, and their relations with all n nodes are computed (hierarchical selection). In case $d_ftf <= ftt$, n is increased, s.t. $d_ftf > ftt$. This might lead to a large number of pairs of nodes to process at a given iteration step. In order to control memory consumption, one might therefore set max_pairs to a suitable value, triggering a subiteration if this value is exceeded.

In order to parallelize the iterative computation, one may pass the arguments from_pos and to_pos. They determine the range of **source nodes** to process (endpoint excluded). Hence, from_pos has to be in [0, g.n[,

and to_pos in [1,g.n]. For instance, given the node table above

>>> g.v time x 0 -3.6 -3 1 -1.1 3 2 1.4 1 3 4.0 12 4 6.3 7

we can compute all relations of the source nodes in [1,3[by

>>> g.create_edges_ft(ft_feature=('time', 5), from_pos=1, to_pos=3)

>>> g.e
 ft_r
s t
1 2 2.5
2 3 2.6
4 4.9

Like create_edges, this method also works with a pd.HDFStore containing the DataFrame representing the node table. Only the data requested by ft_feature, transfer_features and the user-defined connectors and selectors at each iteration step is then pulled from the store. The node table in the store has to be in table(t) format, and additionally, the fast_track feature has to be a data column. For instance, storing the above node table

```
>>> vstore = pd.HDFStore('vstore.h5')
>>> vstore.put('node_table', v, format='t', data_columns=True,
... index=False)
```

one may initiate a DeepGraph instance with the store

>>> g = dg.DeepGraph(vstore)

```
>>> g.v
<class 'pandas.io.pytables.HDFStore'>
File path: vstore.h5
/node_table frame_table (typ->appendable,nrows->5,ncols->2,
indexers->[index],dc->[time,x])
```

and then create edges the same way as if g.v were a DataFrame

```
>>> g.create_edges_ft(ft_feature=('time', 5), from_pos=1, to_pos=3)
```

>>> g.e
 ft_r
s t
1 2 2.5
2 3 2.6
4 4.9

Warning: There is no assertion whether the node table in a store is sorted by the fast-track feature! The result of an unsorted table is unpredictable, and generally not correct.

Parameters

- **ft_feature** (*tuple*) A tuple (ftf, ftt), where ftf is a column name of v (the fast-track feature) and ftt a positive number (the fast-track threshold). The fast-track feature may contain integers or floats, but datetime-like values are also accepted. In that case, ft_feature has to be a tuple of length 3, (ftf, ftt, dt_unit), where dt_unit is on of {'D','h','m','s','ms','us','ns'}:
 - D: days
 - -h: hours
 - *m*: minutes
 - s: seconds
 - *ms*: milliseconds
 - us: microseconds
 - ns: nanoseconds

determining the unit in which the temporal distance is measured. The variable name of the fast-track relation transferred to e is ft_r.

- connectors (function or array_like, optional (default=None)) -User defined connector function(s) that compute pairwise relations between the nodes in v. A connector accepts multiple column names of v (with '_s' and/or '_t' appended, indicating source node values and target node values, respectively) as input, as well as already computed relations of former connectors. A connector function may have multiple output variables. Every output variable has to be a 1-dimensional np.ndarray (with arbitrary dtype, including object). A connector may also depend on the fast- track relations ('ft_r'). See dg.functions for examplary connector functions.
- **selectors** (function or array_like, optional (default=None)) User defined selector function(s) that select edges during the iteration process, based on some conditions on the node's features and their computed relations. Every selector function must have sources and targets as input arguments as well as in the return statement. A selector may depend on column names of v (with '_s' and/or '_t' appended) and/or computed relations of connector functions, and/or computed relations of former selector functions. Apart from sources and targets, they may also return computed relations (see connectors). A selector may also depend on the fast-track relations ('ft_r'). See dg. functions for exemplary selector functions.

Note: To specify the hierarchical order of the selection by the fast-track selector, insert the string 'ft_selector' in the corresponding position of the selectors list. Otherwise, computation of ft_r and selection by the fast-track selector is carried out first.

- transfer_features (str, int or array_like, optional (default=None)) A (list of) column name(s) of v, indicating which features of v to transfer to e (appending '_s' and '_t' to the column names of e, indicating source and target node features, respectively).
- **r_dtype_dic** (*dict*, *optional* (*default=None*)) A dictionary with names of computed relations of connectors and/or selectors as keys and dtypes as values. Forces the data types of the computed relations in e during the iteration (but **after** all selectors and connectors were processed), otherwise infers them.
- **no_transfer_rs** (*str* or *array_like*, *optional* (*default=None*)) Name(s) of computed relations that are not to be transferred to the created edge table e.

Can be used to save memory, e.g., if a selector depends on computed relations that are of no interest otherwise.

- **min_chunk_size** (*int*, *optional* (*default=1000*)) The minimum number of nodes to form pairs of at each iteration step. See above for details.
- **max_pairs** (positive integer, optional (default=1e6)) The maximum number of pairs of nodes to process at any given iteration step. If the number is exceeded, a memory saving subiteration is applied.
- **from_pos** (*int*, *optional* (*default=0*)) The locational index (.iloc) of v to start the iteration. Determines the range of **source nodes** to process, in conjuction with to_pos. Has to be in [0, g.n[, and smaller than to_pos. See above for details and an example.
- to_pos (*int*, *optional* (*default=None*)) The locational index (.iloc) of v to end the iteration (excluded). Determines the range of **source nodes** to process, in conjuction with from_pos. Has to be in [1, g.n], and larger than from_pos. Defaults to None, which translates to the last node of v, to_pos=g.n. See above for details and an example.
- hdf_key (str, optional (default=None)) If you initialized dg. DeepGraph with a pandas.HDFStore and the store has multiple nodes, you must pass the key to the node in the store that corresponds to the node table.
- **verbose**(*bool*, *optional* (*default=False*)) Whether to print information at each step of the iteration process.
- **logfile** (*str*, *optional* (*default=None*)) Create a log-file named by logfile. Contains the time and date of the method's call, the input arguments and time mesaurements for each iteration step. A plot of logfile can be created by dg. DeepGraph.plot_logfile.

Returns e – Set the created edge table e as attribute of dg.DeepGraph.

Return type pd.DataFrame

See also:

create_edges()

Notes

The parameter min_chunk_size enforces a vectorized iteration and changing its value can both accelerate or slow down computation time. This depends mostly on the distribution of values of the fast track feature, and the complexity of the given connectors and selectors. Use the logging capabilites to determine a good value.

When using a pd.HDFStore for the computation, the following advice might be considered. Recall that the only requirements on the node in the store are: the format is table(t), not fixed(t); the node is sorted by the fast-track feature; and the fast-track feature is a data column.

The recommended procedure of storing a given node table v in a store is the following (using the above node table):

```
>>> vstore = pd.HDFStore('vstore.h5')
>>> vstore.put('node_table', v, format='t', data_columns=True,
... index=False)
```

Setting index=False significantly decreases the time to construct the node in the store, and also reduces the resulting file size. It has no impact, however, on the capability of querying the store (with the pd.HDFStore.select* methods).

However, there are two reasons one might want to create a pytables index of the fast-track feature:

1. The node table might be too large to be sorted in memory. To sort it on disc, one may proceed as follows. Assuming an unsorted (large) node table

```
>>> v = pd.DataFrame({'time': [6.3,-3.6,4.,-1.1,1.4],
... 'x': [-3.,3.,1.,12.,7.]})
```

>>> v
 time x
0 6.3 -3
1 -3.6 3
2 4.0 1
3 -1.1 12
4 1.4 7

one stores it as recommended

```
>>> vstore = pd.HDFStore('vstore.h5')
>>> vstore.put('node_table', v, format='t', data_columns=True,
... index=False)
>>> vstore.get_storer('node_table').group.table
/node_table/table (Table(5,)) ''
description := {
    "index": Int64Col(shape=(), dflt=0, pos=0),
    "time": Float64Col(shape=(), dflt=0.0, pos=1),
    "x": Float64Col(shape=(), dflt=0.0, pos=2)}
byteorder := 'little'
chunkshape := (2730,)
```

creates a (full) pytables index of the fast-track feature

and then sorts it on disc with

```
>>> vstore.close()
>>> !ptrepack --chunkshape=auto --sortby=time vstore.h5 s_vstore.h5
>>> s_vstore = pd.HDFStore('s_vstore.h5')
```

```
>>> s_vstore.node_table
   time x
```

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1	-3.6	3
3	-1.1	12
4	1.4	7
2	4.0	1
0	6.3	-3

2. To speed up the internal queries on the fast-track feature

```
>>> s_vstore.create_table_index('node_table', columns=['time'],
... kind='full')
```

See http://stackoverflow.com/questions/17893370/ptrepack-sortby-needs-full-index and https://gist.github. com/michaelaye/810bd0720bb1732067ff for details, benchmarks, and the effects of compressing the store.

Graph Partitioning

<pre>DeepGraph.partition_nodes(features[,])</pre>	Return a supernode DataFrame sv.
<pre>DeepGraph.partition_edges([relations,])</pre>	Return a superedge DataFrame se.
<pre>DeepGraph.partition_graph(features[,])</pre>	Return supergraph DataFrames sv and se.

deepgraph.deepgraph.DeepGraph.partition_nodes

DeepGraph.partition_nodes (*features*, *feature_funcs=None*, *n_nodes=True*, *return_gv=False*) Return a supernode DataFrame sv.

This is essentially wrapper around the pandas groupby method: а SV = v.groupby(features).agg(feature funcs). It creates a (intersection) partition of the nodes in v by the type(s) of feature(s) features, resulting in a supernode DataFrame sv. By passing a dictionary of functions on the features of v, feature_funcs, one may aggregate user-defined values of the partition's elements, the supernodes' features. If n nodes is True, create a column with the number of each supernode's constituent nodes. If return_qv is True, return the created groupby object to facilitate additional operations, such as qv.apply(func, *args, **kwargs).

For details, type help(v.groupby), and/or inspect the available methods of gv.

For examples, see below. For an in-depth description and mathematical details of graph partitioning, see https://arxiv.org/pdf/1604.00971v1.pdf, in particular Sec. III A, E and F.

Parameters

- **features** (*str*, *int or array_like*) Column name(s) of v, indicating the type(s) of feature(s) used to induce a (intersection) partition. Creates a pandas groupby object, gv = v.groupby(features).
- feature_funcs (dict, optional (default=None)) Each key must be a column name of v, each value either a function, or a list of functions, working when passed a pandas.DataFrame or when passed to pandas.DataFrame.apply. See the docstring of gv.agg for details: help(gv.agg).
- **n_nodes** (*bool*, *optional* (*default=True*)) Whether to create a n_nodes column in sv, indicating the number of nodes in each supernode.
- **return_gv** (bool, optional (default=False)) If True, also return the v.groupby(features) object, gv.

Returns

- sv (pd.DataFrame) The aggreated DataFrame of supernodes, sv.
- **gv** (*pandas.core.groupby.DataFrameGroupBy*) The pandas groupby object, v.groupby(features).

See also:

```
partition_edges(), partition_graph()
```

Notes

Currently, NA groups in GroupBy are automatically excluded (silently). One workaround is to use a placeholder (e.g., -1, 'none') for NA values before doing the groupby (calling this method). See http://stackoverflow.com/ questions/18429491/groupby-columns-with-nan-missing-values and https://github.com/pydata/pandas/issues/ 3729.

Examples

First, we need a node table, in order to demonstrate its partitioning:

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'x': [-3.4,2.1,-1.1,0.9,2.3],
                      'time': [0,0,2,2,9],
. . .
                      'color': ['g','g','b','g','r'],
. . .
                      'size': [1,3,2,3,1]})
. . .
>>> g = dg.DeepGraph(v)
>>> g.v
 color size time
                       Х
0
      g
         1
              0 -3.4
           3
                 0 2.1
1
      g
2
           2
                 2 -1.1
      b
3
            3
                  2 0.9
      g
                  9 2.3
4
           1
      r
```

Create a partition by the type of feature 'color':

Create an intersection partition by the types of features 'color' and 'size' (which is a further refinement of the last partition):

Partition by 'color' and collect x values:

Partition by 'color' and aggregate with different functions:

```
>>> g.partition_nodes('color', {'time': [lambda x: list(x), np.max],
                                 'x': [np.mean, np.sum, np.std]})
. . .
                                     x_std time_<lambda> time_amax
       n_nodes
                   x_mean
                          x_sum
color
                                                                    2
b
             1 -1.100000
                            -1.1
                                       NaN
                                                      [2]
                            -0.4 2.891943
                                                [0, 0, 2]
                                                                    2
g
             3 -0.133333
                                                                    9
r
             1
                2.300000
                             2.3
                                       NaN
                                                      [9]
```

deepgraph.deepgraph.DeepGraph.partition_edges

DeepGraph.partition_edges (relations=None, source_features=None, target_features=None, relation_funcs=None, n_edges=True, return_ge=False)

Return a superedge DataFrame se.

This method allows you to partition the edges in e by their types of relations, but also by the types of features of their incident source and target nodes, and any combination of the three.

Essentially, this method is a wrapper around the pandas groupby method: se = e.groupby(relations + features_s + features_t).agg(relation_funcs), where relations are column names of e, and in order to group e by features_s and/or features_t, the features of type source_features and/or target_features (column names of v) are transferred to e, appending '_s' and/or '_t' to the corresponding column names of e (if they are not already present). The only requirement on the combination of relations, source_features and target_features is that at least on of the lists has to be of length >= 1.

By passing a dictionary of functions on the relations of e, relation_funcs, one may aggregate user-defined values of the partition's elements, the superedges' relations. If n_edges is True, create a column with the number of each superedge's constituent edges. If return_ge is True, return the created groupby object to facilitate additional operations, such as ge.apply(func, *args, **kwargs).

For details, type help(g.e.groupby), and/or inspect the available methods of ge.

For examples, see below. For an in-depth description and mathematical details of graph partitioning, see https: //arxiv.org/pdf/1604.00971v1.pdf, in particular Sec. III B, E and F.

Parameters

- relations (str, int or array_like, optional (default=None)) Column name(s) of e, indicating the type(s) of relation(s) used to induce a (intersection) partition of e (in conjunction with source_features and target_features).
- source_features (str, int or array_like, optional (default=None)) Column name(s) of v, indicating the type(s) of feature(s) of the edges' incident source nodes used to induce a (intersection) partition of e (in conjunction with relations and target_features).
- target_features (str, int or array_like, optional (default=None)) Column name(s) of v, indicating the type(s) of feature(s) of

the edges' incident target nodes used to induce a (intersection) partition of e (in conjunction with relations and source_features).

- relation_funcs (dict, optional (default=None)) Each key must be a column name of e, each value a (list of) function(s), working when passed a pandas. DataFrame or when passed to pandas.DataFrame.apply. See the docstring of ge.agg for details: help(ge.agg).
- **n_edges** (*bool*, *optional* (*default=True*)) Whether to create a n_edges column in se, indicating the number of edges in each superedge.
- **return_ge** (*bool*, *optional* (*default=False*)) If True, also return the pandas groupby object, ge.

Returns

- se (pd.DataFrame) The aggreated DataFrame of superedges, se.
- ge (pandas.core.groupby.DataFrameGroupBy) The pandas groupby object, ge.

See also:

```
partition_nodes(), partition_graph()
```

Notes

Currently, NA groups in GroupBy are automatically excluded (silently). One workaround is to use a placeholder (e.g., -1, 'none') for NA values before doing the groupby (calling this method). See http://stackoverflow.com/ questions/18429491/groupby-columns-with-nan-missing-values and https://github.com/pydata/pandas/issues/ 3729.

Examples

First, we need to create a graph in order to demonstrate how to partition its edge set.

Create a node table:

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'x': [-3.4,2.1,-1.1,0.9,2.3],
... 'time': [0,1,2,5,9],
... 'color': ['g','g','b','g','r'],
... 'size': [1,3,2,3,1]})
>>> g = dg.DeepGraph(v)
```

```
>>> g.v
 color size time
               Х
   g 1 0-3.4
0
        3
            1 2.1
1
    q
            2 -1.1
2
    b
        2
3
        3
            5 0.9
    g
4
        1
            9 2.3
    r
```

Create an edge table:

```
>>> def some_relations(ft_r, x_s,x_t,color_s,color_t,size_s,size_t):
... dx = x_t - x_s
... v = dx / ft_r
```

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```
... same_color = color_s == color_t
... larger_than = size_s > size_t
... return dx, v, same_color, larger_than
>>> g.create_edges_ft(('time', 5), connectors=some_relations)
>>> g.e.rename(columns={'ft_r': 'dt'}, inplace=True)
>>> g.e['inds'] = g.e.index.values # to ease the eyes
```

>>	->	g.e						
		dx	dt	larger_than	same_color	V	ir	nds
S	t							
0	1	5.5	1	False	True	5.500000	(0,	1)
	2	2.3	2	False	False	1.150000	(0,	2)
	3	4.3	5	False	True	0.860000	(0,	3)
1	2	-3.2	1	True	False	-3.200000	(1,	2)
	3	-1.2	4	False	True	-0.300000	(1,	3)
2	3	2.0	3	False	False	0.666667	(2,	3)
3	4	1.4	4	True	False	0.350000	(3,	4)

Partitioning by the type of relation 'larger_than':

>>> g.partit	<pre>>>> g.partition_edges(relations='larger_than',</pre>											
		relation_funcs=	{'dx': ['	<pre>mean', 'std'],</pre>								
			'same_co	<pre>lor': 'sum'})</pre>								
	n_edges	same_color_sum	dx_mean	dx_std								
larger_than												
False	5	3	2.58	2.558711								
True	2	0	-0.90	3.252691								

A refinement of the last partition by the type of relation 'same_color':

```
>>> g.partition_edges(relations=['larger_than', 'same_color'],
                     relation_funcs={'dx': ['mean', 'std'],
• • •
                                     'dt': lambda x: tuple(x) })
. . .
                       n_edges dt_<lambda> dx_mean dx_std
larger_than same_color
False
      False
                             2
                                   (2, 3) 2.150000 0.212132
           True
                             3 (1, 5, 4) 2.866667 3.572581
True
           False
                             2
                                   (1, 4) - 0.900000 3.252691
```

Partitioning by the type of source feature 'color':

As one can see, the type of feature 'color' of the source nodes has been transferred to e:

>>>	g.e							
	dx	dt	larger_than	same_color	V	inds	color_s	
st								
0 1	5.5	1	False	True	5.500000	(0, 1)	g	
2	2.3	2	False	False	1.150000	(0, 2)	g	
3	4.3	5	False	True	0.860000	(0, 3)	g	

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1 2 -3.2	1	True	False -3.200000	(1, 2)	g	
3 -1.2	4	False	True -0.300000	(1, 3)	g	
2 3 2.0	3	False	False 0.666667	(2, 3)	b	
3 4 1.4	4	True	False 0.350000	(3, 4)	g	

A further refinement of the last partition by the type of source feature 'size':

>>> g	g.partitior	n_edges(son	urce_feature	s=['color', 'size'],
		re	lation_funcs	={'same_color': 'sum',
				<pre>'inds': lambda x: tuple(x)})</pre>
		n_edges	same_color	inds
color	_s size_s			
b	2	1	0	((2, 3),)
g	1	3	2	((0, 1), (0, 2), (0, 3))
	3	3	1	((1, 2), (1, 3), (3, 4))

Partitioning by the types of target features ('color', 'size'):

>>> g.p	<pre>>>> g.partition_edges(target_features=['color', 'size'],</pre>										
	<pre> relation_funcs={'same_color': 'sum',</pre>										
				<pre>'inds': lambda x: tuple(x) })</pre>							
		n_edges	same_color	inds							
color_t	size_t										
b	2	2	0	((0, 2), (1, 2))							
g	3	4	3	((0, 1), (0, 3), (1, 3), (2, 3))							
r	1	1	0	((3, 4),)							

Partitioning by the type of source feature 'color' and the type of target feature 'size':

>>> g.	<pre>>>> g.partition_edges(source_features='color', target_features='size',</pre>							
		re	lation_funcs	={'same_color': 'sum',				
	'inds': lambda x: tuple(x)})							
		n_edges	same_color	inds				
color_	_s size_t							
b	3	1	0	((2, 3),)				
g	1	1	0	((3, 4),)				
	2	2	0	((0, 2), (1, 2))				
	3	3	3	((0, 1), (0, 3), (1, 3))				

A further refinement of the last partition by the type of relation 'larger_than':

<pre>>>> g.partition_edges(relations='larger_than',</pre>							
		source	e_features='c	color', target_features='size',			
		relati	on_funcs={'i	<pre>inds': lambda x: tuple(x) })</pre>			
			n_edges	inds			
larger_than	color_s	size_t					
False	b	3	1	((2, 3),)			
	g	2	1	((0, 2),)			
		3	3 ((0	0, 1), (0, 3), (1, 3))			
True	g	1	1	((3, 4),)			
		2	1	((1, 2),)			

deepgraph.deepgraph.DeepGraph.partition_graph

Return supergraph DataFrames sv and se.

This method allows partitioning of the graph represented by v and e into a supergraph, sv and se. It creates a (intersection) partition of the nodes in v by the type(s) of feature(s) features, together with the (intersection) partition's **corresponding** partition of the edges in e.

Essentially, this method is а wrapper around pandas methgroupby ods: v.groupby(features).agg(feature funcs) and se sv = e.groupby(features_s+features_t).agg(relation_funcs). In order to group e by features_s and features t, the features of type features are transferred to e, appending 's' and 't' to the corresponding column names of e, indicating source and target features, respectively (if they are not already present).

By passing a dictionary of functions on the features (relations) of v (e), feature_funcs (relation_funcs), one may aggregate user-defined values of the partition's elements, the supernodes' (superedges') features (relations). If n_nodes (n_edges) is True, create a column with the number of each supernode's (superedge's) constituent nodes (edges).

If return_gve is True, return the created groupby objects to facilitate additional operations, such as gv.apply(func, *args, **kwargs) or ge.apply(func, *args, **kwargs).

For details, type help(g.v.groupby), and/or inspect the available methods of gv.

For examples, see below. For an in-depth description and mathematical details of graph partitioning, see https://arxiv.org/pdf/1604.00971v1.pdf, in particular Sec. III C, E and F.

Parameters

- **features** (*str*, *int or array_like*) Column name(s) of v, indicating the type(s) of feature(s) used to induce a (intersection) partition of v, and its **corresponding** partition of the edges in e. Creates pandas groupby objects, gv and ge.
- feature_funcs (dict, optional (default=None)) Each key must be a column name of v, each value either a function, or a list of functions, working when passed a pandas.DataFrame or when passed to pandas.DataFrame.apply. See the docstring of gv.agg for details: help(gv.agg).
- relation_funcs (dict, optional (default=None)) Each key must be a column name of e, each value either a function, or a list of functions, working when passed a pandas.DataFrame or when passed to pandas.DataFrame.apply. See the doc-string of ge.agg for details: help(ge.agg).
- **n_nodes** (*bool*, *optional* (*default=True*)) Whether to create a n_nodes column in sv, indicating the number of nodes in each supernode.
- **n_edges** (*bool*, *optional* (*default=True*)) Whether to create a n_edges column in se, indicating the number of edges in each superedge.
- **return_gve**(bool, optional (default=False))-If True, also return the pandas groupby objects, gv and ge.

Returns

- sv (pd.DataFrame) The aggreated DataFrame of supernodes, sv.
- se (pd.DataFrame) The aggregated DataFrame of superedges, se.
- **gv** (*pandas.core.groupby.DataFrameGroupBy*) The pandas groupby object, v.groupby(features).

• **ge** (*pandas.core.groupby.DataFrameGroupBy*) – The pandas groupby object, e.groupby(features_i+feaures_j).

See also:

```
partition_nodes(), partition_edges()
```

Notes

Currently, NA groups in GroupBy are automatically excluded (silently). One workaround is to use a placeholder (e.g., -1, 'none') for NA values before doing the groupby (calling this method). See http://stackoverflow.com/ questions/18429491/groupby-columns-with-nan-missing-values and https://github.com/pydata/pandas/issues/ 3729.

Examples

First, we need to create a graph in order to demonstrate its partitioning into a supergraph.

Create a node table:

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'x': [-3.4,2.1,-1.1,0.9,2.3],
... 'time': [0,1,2,5,9],
... 'color': ['g','g','b','g','r'],
... 'size': [1,3,2,3,1]})
>>> g = dg.DeepGraph(v)
```

>>> g.v

	-			
	color	size	time	Х
0	g	1	0	-3.4
1	g	3	1	2.1
2	b	2	2	-1.1
3	g	3	5	0.9
4	r	1	9	2.3

Create an edge table:

```
>>> def some_relations(ft_r, x_s,x_t,color_s,color_t,size_s,size_t):
... dx = x_t - x_s
... v = dx / ft_r
... same_color = color_s == color_t
... larger_than = size_s > size_t
... return dx, v, same_color, larger_than
>>> g.create_edges_ft(('time', 5), connectors=some_relations)
>>> g.e.rename(columns={'ft_r': 'dt'}, inplace=True)
>>> g.e['inds'] = g.e.index.values # to ease the eyes
```

>	>>	g.e						
		dx	dt	larger_than	same_color	V	in	ds
s	t							
0	1	5.5	1	False	True	5.500000	(0,	1)
	2	2.3	2	False	False	1.150000	(0,	2)
	3	4.3	5	False	True	0.860000	(0,	3)
1	2	-3.2	1	True	False	-3.200000	(1,	2)
	3	-1.2	4	False	True	-0.300000	(1,	3)

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23	2.0	3	False	False	0.666667	(2,	3)
34	1.4	4	True	False	0.350000	(3,	4)

Create a supergraph by partitioning by the type of feature 'color':

```
>>> sv, se = g.partition_graph('color')
```

>>> sv	
	n_nodes
color	
b	1
g	3
r	1

>>> se

```
n_edges
color_s color_t
b g 1
g b 2
g 3
r 1
```

Create intersection partitions by the types of features 'color' and 'size' (which are further refinements of the last partitions):

```
>>> sv, se = g.partition_graph(
       ['color', 'size'],
       relation_funcs={'inds': lambda x: tuple(x)})
```

```
>>> sv
```

```
n_nodes
color size
b 2 1
g 1 1
3 2
r 1 1
```

>>> se							
				n_edges		inds	
color_s	size_s	color_t	size_t				
b	2	g	3	1	((2,	3),)	
g	1	b	2	1	((0,	2),)	
		g	3	2	((0, 1), (0,	3))	
	3	b	2	1	((1,	2),)	
		g	3	1	((1,	3),)	
		r	1	1	((3,	4),)	

Partition by 'color' and aggregate some properties:

time

```
>>> sv, se = g.partition_graph('color',
... feature_funcs={'time': lambda x: list(x)},
... relation_funcs={'larger_than': 'sum', 'same_color': 'sum'})
```

```
>>> sv
n_nodes
```

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color					
b	1	[2]			
g	3	[0, 1, 5]			
r	1	[9]			
>>> se					
		n_edges lar	ger_than	same_color	
color_s	color_t				
b	g	1	False	0	
g	b	2	True	0	
	g	3	False	3	
	r	1	True	0	

Graph Interfaces

DeepGraph.return_cs_graph([relations,	Return scipy.sparse.coo_matrix representa-
dropna])	tion(s).
<pre>DeepGraph.return_nx_graph([features,])</pre>	Return a networkx.DiGraph representation.
DeepGraph.return_nx_multigraph([features,	Return a networkx.MultiDiGraph representa-
])	tion.
<pre>DeepGraph.return_gt_graph([features,])</pre>	Return a graph_tool.Graph representation.

deepgraph.deepgraph.DeepGraph.return_cs_graph

DeepGraph.return_cs_graph(relations=False, dropna=True)

Return scipy.sparse.coo_matrix representation(s).

Create a compressed sparse graph representation for each type of relation given by relations. relations can either be False, True, or a (list of) column name(s) of e. If relations is False (default), return a single csgraph entailing all edges in e.index, each with a weight of 1 (in that case, dropna is discarded). If relations is True, create one csgraph for each column of e, where the weights are given by the columns' values. If only a subset of columns is to be mapped to csgraphs, relations has to be a (list of) column name(s) of e.

The argument dropna indicates whether to discard edges with NA values or not. If dropna is True or False, it applies to all types of relations given by relations. However, dropna can also be array_like with the same shape as relations (or with the same shape as e.columns, if relations is True).

Parameters

- **relations** (bool, str or array_like, optional (default=False)) The types of relations to be mapped to scipy csgraphs. Can be False, True, or a (list of) column name(s) of e.
- dropna (bool or array_like, optional (default=True)) Whether to drop edges with NA values. If True or False, applies to all relations given by relations. Otherwise, must be the same shape as relations. If relations is False, dropna is discarded.
- **Returns csgraph** A dictionary, where keys are column names of e, and values are the corresponding scipy.sparse.coo_matrix instance(s). If only one csgraph is created, return it directly.

Return type scipy.sparse.coo_matrix or dict

See also:

return_nx_graph(), return_nx_multigraph(), return_gt_graph()

deepgraph.deepgraph.DeepGraph.return_nx_graph

DeepGraph.return_nx_graph (features=False, relations=False, dropna='none')

Return a networkx.DiGraph representation.

Create a networkx.DiGraph representation of the graph given by v and e. Node and edge properties to transfer can be indicated by the features and relations input arguments. Whether to drop edges with NA values in the subset of types of relations given by relations can be controlled by dropna.

Needs pandas $\geq 0.17.0$.

Parameters

- **features** (bool, str, or array_like, optional (default=False)) Indicates which types of features to transfer as node attributes. Can be column name(s) of v, False or True. If False, create no node attributes. If True, create node attributes for every column in v. If str or array_like, must be column name(s) of v indicating which types of features to transfer.
- **relations** (bool, str, or array_like, optional (default=False)) - Indicates which types of relations to transfer as edge attributes. Can be column name(s) of e, False or True. If False, create no edge attributes (all edges in e.index are transferred, regardless of dropna). If True, create edge attributes for every column in e (all edges in e.index are transferred, regardless of dropna). If str or array_like, must be column name(s) of e indicating which types of relations to transfer (which edges are transferred can be controlled by dropna).
- **dropna** (*str*, *optional* (*default='none'*)) One of {'none', 'any', 'all'}. If 'none', all edges in e.index are transferred. If 'any', drop all edges (rows) in e[relations] where any NA values are present. If 'all', drop all edges (rows) in e[relations] where all values are NA. Only has an effect if relations is str or array_like.

Returns nx_g

Return type networkx.DiGraph

See also:

return_nx_multigraph(), return_cs_graph(), return_gt_graph()

deepgraph.deepgraph.DeepGraph.return_nx_multigraph

DeepGraph.return_nx_multigraph (*features=False*, *relations=False*, *dropna=True*) Return a networkx.MultiDiGraph representation.

Create a networkx.MultiDiGraph representation of the graph given by v and e. As opposed to return_nx_graph, where every row of e is treated as one edge, this method treats every cell of e as one edge. The input argument features indicates which node properties to transfer. relations indicates which edges to transfer. Whether to drop edges with NA values can be controlled by dropna.

Needs pandas >= 0.17.0.

Parameters

- **features** (bool, str, or array_like, optional (default=False)) Indicates which types of features to transfer as node attributes. Can be column name(s) of v, False or True. If False, create no node attributes. If True, create node attributes for every column in v. If str or array_like, must be column name(s) of v indicating which types of features to transfer.
- relations (bool, str, or array_like, optional (default=False)) - Indicates which cells of e to transfer as edges. Can be False, True, or a (list of) column name(s) of e. If False (default), all cells of e are translated to edges, but their values are not transferred as edge attributes. If True, all cells of e are translated, and their values are transferred as edge attributes. If str or array_like, must be column name(s) of e, restricting the translation of cells to edges to e[relations] (values are transferred as edge attributes).
- **dropna** (bool, optional (default=True)) Whether to drop edges with NA values. Cells in e with NA values are not translated to edges.

Returns nx_g

Return type networkx.MultiDiGraph

See also:

return_nx_graph(), return_cs_graph(), return_gt_graph()

deepgraph.deepgraph.DeepGraph.return_gt_graph

DeepGraph.return_gt_graph (features=False, relations=False, dropna='none', node_indices=False, edge indices=False)

Return a graph_tool.Graph representation.

Create a graph_tool.Graph (directed) representation of the graph given by v and e. Node and edge properties to transfer can be indicated by the features and relations input arguments. Whether to drop edges with NA values in the subset of types of relations given by relations can be controlled by dropna. If the nodes in v are not indexed by consecutive integers starting from 0, one may internalize the original node and edge indices as propertymaps by setting node_indices and/or edge_indices to True.

Parameters

- **features** (bool, str, or array_like, optional (default=False))-Indicates which types of features to internalize as graph_tool.PropertyMap. Can be column name(s) of v, False or True. If False, create no propertymaps. If True, create propertymaps for every column in v. If str or array_like, must be column name(s) of v indicating which types of features to internalize.
- relations (bool, str, or array_like, optional (default=False)) - Indicates which types of relations to internalize as graph_tool.PropertyMap. Can be column name(s) of e, False or True. If False, create no propertymaps (all edges in e.index are transferred, regardless of dropna). If True, create propertymaps for every column in e (all edges in e.index are transferred, regardless of dropna). If str or array_like, must be column name(s) of e indicating which types of relations to internalize (which edges are transferred can be controlled by dropna).
- **dropna** (*str*, *optional* (*default='none'*)) One of {'none', 'any', 'all'}. If 'none', all edges in e.index are transferred. If 'any', drop all edges (rows) in e[relations] where any NA values are present. If 'all', drop all edges (rows) in e[relations] where all values are NA. Only has an effect if relations is str or array_like.

- **node_indices** (*bool*, *optional* (*default=False*)) If True, internalize a vertex propertymap i with the original node indices.
- **edge_indices** (*bool*, *optional* (*default=False*)) If True, internalize edge propertymaps s and t with the original source and target node indices of the edges, respectively.

Returns gt_g

Return type graph_tool.Graph

See also:

return_cs_graph(), return_nx_graph(), return_nx_multigraph()

Notes

If the index of v is not pd.RangeIndex(start=0,stop=len(v), step=1), the indices will be enumerated, which is expensive for large graphs.

Plotting Methods

<pre>DeepGraph.plot_2d(x,y[,edges,C,])</pre>	Plot nodes and corresponding edges in 2 dimensions.
<pre>DeepGraph.plot_2d_generator(x, y, by[,])</pre>	Plot nodes and corresponding edges by groups.
<pre>DeepGraph.plot_map(lon, lat[, edges, C,])</pre>	Plot nodes and corresponding edges on a basemap.
<pre>DeepGraph.plot_map_generator(lon, lat, by[,</pre>	Plot nodes and corresponding edges by groups, on
])	basemaps.
<pre>DeepGraph.plot_hist(x[, bins, log_bins,])</pre>	Plot a histogram (or pdf) of x.
<pre>DeepGraph.plot_logfile(logfile)</pre>	Plot a logfile.

deepgraph.deepgraph.DeepGraph.plot_2d

Plot nodes and corresponding edges in 2 dimensions.

Create a scatter plot of the nodes in v, and optionally a quiver plot of the corresponding edges in e.

The xy-coordinates of the scatter plot are determined by the values of v[x] and v[y], where x and y are column names of v (the arrow's coordinates are determined automatically).

In order to map colors to the arrows, either C or C_split_0 can be be passed, an array of the same length as e. Passing C creates a single quiver plot (qu). Passing C_split_0 creates two separate quiver plots, one for all edges where C_split_0 == 0 (qu_0), and one for all other edges (qu). By default, the arrows of qu_0 have no head, indicating "undirected" edges. This can be useful, for instance, when C_split_0 represents an array of temporal distances.

In order to control the plotting parameters of the scatter, quiver and/or quiver_0 plots, one may pass keyword arguments by setting kwds_scatter, kwds_quiver and/or kwds_quiver_0.

Can be used iteratively by passing ax.

Parameters

• \mathbf{x} (*int* or *str*) – A column name of v, determining the x-coordinates of the scatter plot of nodes.

- **y** (*int* or *str*) A column name of v, determining the y-coordinates of the scatter plot of nodes.
- edges (bool, optional (default=True)) Whether to create a quiver plot (2-D field of arrows) of the edges between the nodes.
- C (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. Has no effect if C_split_0 is passed as an argument.
- C_split_0 (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. If this parameter is passed, C has no effect, and two separate quiver plots are created (qu and qu_0).
- **kwds_scatter** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to scatter.
- kwds_quiver (dict, optional (default=None)) kwargs to be passed to quiver (qu).
- kwds_quiver_0 (dict, optional (default=None)) kwargs to be passed to quiver (qu_0). Only has an effect if C_split_0 has been set.
- **ax** (matplotlib axes object, optional (default=None)) An axes instance to use.
- **Returns obj** If C_split_0 has been passed, return a dict of matplotlib objects with the following keys: ['fig', 'ax', 'pc', 'qu', 'qu_0']. Otherwise, return a dict with keys: ['fig', 'ax', 'pc', 'qu'].

Return type dict

Notes

When passing C_split_0, the color of the arrows in qu_0 can be set by passing the keyword argument *color* to kwds_quiver_0. The color of the arrows in qu, however, are determined by C_split_0.

The default drawing order is set to: 1. quiver_0 (zorder=1) 2. quiver (zorder=2) 3. scatter (zorder=3) This order can be changed by setting the zorder in kwds_quiver_0, kwds_quiver and/or kwds_scatter. See also http://matplotlib.org/examples/pylab_examples/zorder_demo.html

See also:

plot_2d_generator(), plot_3d(), plot_map(), plot_map_generator()

deepgraph.deepgraph.DeepGraph.plot_2d_generator

DeepGraph.plot_2d_generator(x, y, by, edges=False, C=None, C_split_0=None, kwds_scatter=None, kwds_quiver=None, kwds_quiver_0=None, passable_ax=False)

Plot nodes and corresponding edges by groups.

Create a generator of scatter plots of the nodes in v, split in groups by v.groupby(by). If edges is set True, also create a quiver plot of each group's corresponding edges.

The xy-coordinates of the scatter plots are determined by the values of v[x] and v[y], where x and y are column names of v (the arrow's coordinates are determined automatically).

In order to map colors to the arrows, either C or C_split_0 can be be passed, an array of the same length as e. Passing C creates a single quiver plot (qu). Passing C_split_0 creates two separate quiver plots, one for all edges where C_split_0 == 0 (qu_0), and one for all other edges (qu). By default, the arrows of qu_0 have

no head, indicating "undirected" edges. This can be useful, for instance, when C_split_0 represents an array of temporal distances.

When mapping colors to arrows by setting C (or C_split_0), *clim* is automatically set to the min and max values of the entire array. In case one wants clim to be set to min and max values for each group's colors, one may explicitly pass *clim* = None to kwds_quiver.

The same behaviour occurs when passing a sequence of g.n Numbers as colors c to kwds_scatter. In that case, *vmin* and *vmax* are automatically set to c.min() and c.max() of all nodes. Explicitly setting *vmin* and *vmax* to *None*, the min and max values of the groups' color arrays are used.

In order to control the plotting parameters of the scatter, quiver and/or quiver_0 plots, one may pass keyword arguments by setting kwds_scatter, kwds_quiver and/or kwds_quiver_0.

If passable_ax is True, create a generator of functions. Each function takes a matplotlib axes object as input, and returns a scatter/quiver plot.

Parameters

- \mathbf{x} (int or str) A column name of v, determining the x-coordinates of the scatter plot of nodes.
- **y** (*int* or *str*) A column name of v, determining the y-coordinates of the scatter plot of nodes.
- by (array_like) Column name(s) of v, determining the groups to create plots of.
- edges (bool, optional (default=True)) Whether to create a quiver plot (2-D field of arrows) of the edges between the nodes.
- C (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. Has no effect if C_split_0 is passed as an argument.
- C_split_0 (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. If this parameter is passed, C has no effect, and two separate quiver plots are created (qu and qu_0).
- **kwds_scatter** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to scatter.
- **kwds_quiver** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to quiver (qu).
- kwds_quiver_0 (dict, optional (default=None)) kwargs to be passed to quiver (qu_0). Only has an effect if C_split_0 has been set.
- **passable_ax**(bool, optional (default=False))-If True, return a generator of functions. Each function takes a matplotlib axes object as input, and returns a dict of matplotlib objects.
- **Returns obj** If C_split_0 has been passed, return a generator of dicts of matplotlib objects with the following keys: ['fig', 'ax', 'pc', 'qu', 'qu_0', 'group']. Otherwise, return a generator of dicts with keys: ['fig', 'ax', 'pc', 'qu', 'group']. If passable_ax is True, return a generator of functions. Each function takes a matplotlib axes object as input, and returns a dict as described above.

Return type generator

Notes

When passing C_split_0, the color of the arrows in qu_0 can be set by passing the keyword argument *color* to kwds_quiver_0. The color of the arrows in qu, however, are determined by C_split_0.

The default drawing order is set to: 1. quiver_0 (zorder=1) 2. quiver (zorder=2) 3. scatter (zorder=3) This order can be changed by setting the zorder in kwds_quiver_0, kwds_quiver and/or kwds_scatter. See also http://matplotlib.org/examples/pylab_examples/zorder_demo.html

See also:

```
append_binning_labels_v(), plot_2d(), plot_3d(), plot_map(),
plot_map_generator()
```

deepgraph.deepgraph.DeepGraph.plot_map

DeepGraph.plot_map(lon, lat, edges=False, C=None, C_split_0=None, kwds_basemap=None, kwds_scatter=None, kwds_quiver=None, kwds_quiver_0=None, ax=None, m=None)

Plot nodes and corresponding edges on a basemap.

Create a scatter plot of the nodes in v and optionally a quiver plot of the corresponding edges in e on a mpl_toolkits.basemap.Basemap instance.

The coordinates of the scatter plot are determined by the node's longitudes and latitudes (in degrees): v[lon] and v[lat], where lon and lat are column names of v (the arrow's coordinates are determined automatically).

In order to map colors to the arrows, either C or C_split_0 can be be passed, an array of the same length as e. Passing C creates a single quiver plot (qu). Passing C_split_0 creates two separate quiver plots, one for all edges where C_split_0 == 0 (qu_0), and one for all other edges (qu). By default, the arrows of qu_0 have no head, indicating "undirected" edges. This can be useful, for instance, when C_split_0 represents an array of temporal distances.

In order to control the parameters of the basemap, scatter, quiver and/or quiver_0 plots, one may pass keyword arguments by setting kwds_basemap, kwds_scatter, kwds_quiver and/or kwds_quiver_0.

Can be used iteratively by passing ax and/or m.

Parameters

- **lon** (*int* or *str*) A column name of v. The corresponding values must be longitudes in degrees.
- **lat** (*int* or *str*) A column name of v. The corresponding values must be latitudes in degrees.
- edges (bool, optional (default=True)) Whether to create a quiver plot (2-D field of arrows) of the edges between the nodes.
- C (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. Has no effect if C_split_0 is passed as an argument.
- C_split_0 (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. If this parameter is passed, C has no effect, and two separate quiver plots are created (qu and qu_0).
- **kwds_basemap** (*dict*, *optional* (*default=None*)) **kwargs** passed to basemap.

- **kwds_scatter** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to scatter.
- **kwds_quiver** (dict, optional (default=None)) kwargs to be passed to quiver (qu).
- kwds_quiver_0 (dict, optional (default=None)) kwargs to be passed to quiver (qu_0). Only has an effect if C_split_0 has been set.
- **ax** (matplotlib axes object, optional (default=None)) An axes instance to use.
- m (Basemap object, optional (default=None)) A mpl_toolkits.basemap.Basemap instance to use.
- **Returns obj** If C_split_0 has been passed, return a dict of matplotlib objects with the following keys: ['fig', 'ax', 'm', 'pc', 'qu', 'qu_0']. Otherwise, return a dict with keys: ['fig', 'ax', 'm', 'pc', 'qu'].

Return type dict

Notes

When passing C_split_0, the color of the arrows in qu_0 can be set by passing the keyword argument *color* to kwds_quiver_0. The color of the arrows in qu, however, are determined by C_split_0.

The default drawing order is set to: 1. quiver_0 (zorder=1) 2. quiver (zorder=2) 3. scatter (zorder=3) This order can be changed by setting the zorder in kwds_quiver_0, kwds_quiver and/or kwds_scatter. See also http://matplotlib.org/examples/pylab_examples/zorder_demo.html

See also:

plot_map_generator(), plot_2d(), plot_2d_generator(), plot_3d()

deepgraph.deepgraph.DeepGraph.plot_map_generator

DeepGraph.plot_map_generator(lon, lat, by, edges=False, C=None, C_split_0=None, kwds_basemap=None, kwds_scatter=None, kwds_quiver=None, kwds_quiver_0=None, passable_ax=False) Plot nodes and corresponding edges by groups, on basemaps.

Create a generator of scatter plots of the nodes in v, split in groups by v.groupby(by), on a mpl_toolkits. basemap.Basemap instance. If edges is set True, also create a quiver plot of each group's corresponding edges.

The coordinates of the scatter plots are determined by the node's longitudes and latitudes (in degrees): v[lon] and v[lat], where lon and lat are column names of v (the arrow's coordinates are determined automatically).

In order to map colors to the arrows, either C or C_split_0 can be be passed, an array of the same length as e. Passing C creates a single quiver plot (qu). Passing C_split_0 creates two separate quiver plots, one for all edges where C_split_0 == 0 (qu_0), and one for all other edges (qu). By default, the arrows of qu_0 have no head, indicating "undirected" edges. This can be useful, for instance, when C_split_0 represents an array of temporal distances.

When mapping colors to arrows by setting C (or C_split_0), *clim* is automatically set to the min and max values of the entire array. In case one wants clim to be set to min and max values for each group's colors, one may explicitly pass *clim* = None to kwds_quiver.

The same behaviour occurs when passing a sequence of g.n Numbers as colors c to kwds_scatter. In that case, *vmin* and *vmax* are automatically set to c.min() and c.max() of all nodes. Explicitly setting *vmin* and *vmax* to *None*, the min and max values of the groups' color arrays are used.

In order to control the parameters of the basemap, scatter, quiver and/or quiver_0 plots, one may pass keyword arguments by setting kwds_basemap, kwds_scatter, kwds_quiver and/or kwds_quiver_0.

If passable_ax is True, create a generator of functions. Each function takes a matplotlib axes object (and/or a Basemap object) as input, and returns a scatter/quiver plot.

Parameters

- lon (*int* or str) A column name of v. The corresponding values must be longitudes in degrees.
- **lat** (*int* or *str*) A column name of v. The corresponding values must be latitudes in degrees.
- by (array_like) Column name(s) of v, determining the groups to create plots of.
- edges (bool, optional (default=True)) Whether to create a quiver plot (2-D field of arrows) of the edges between the nodes.
- C (array_like, optional (default=None)) An optional array used to map colors to the arrows. Must have the same length es e. Has no effect if C_split_0 is passed as an argument.
- **C_split_0** (*array_like*, *optional* (*default=None*)) An optional array used to map colors to the arrows. Must have the same length es e. If this parameter is passed, C has no effect, and two separate quiver plots are created (qu and qu_0).
- kwds_basemap (dict, optional (default=None)) kwargs passed to basemap.
- **kwds_scatter** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to scatter.
- **kwds_quiver** (*dict*, *optional* (*default=None*)) **kwargs** to be passed to quiver (qu).
- kwds_quiver_0 (dict, optional (default=None)) kwargs to be passed to quiver (qu_0). Only has an effect if C_split_0 has been set.
- **passable_ax**(bool, optional (default=False))-If True, return a generator of functions. Each function takes a matplotlib axes object (and/or a Basemap object) as input, and returns a dict of matplotlib objects.
- **Returns obj**-If C_split_0 has been passed, return a generator of dicts of matplotlib objects with the following keys: ['fig', 'ax', 'm', 'pc', 'qu', 'qu_0', 'group']. Otherwise, return a generator of dicts with keys: ['fig', 'ax', 'm', 'pc', 'qu', 'group']. If passable_ax is True, return a generator of functions. Each function takes a matplotlib axes object (and/or a Basemap object) as input, and returns a dict as described above.

Return type generator

Notes

When passing C_split_0, the color of the arrows in qu_0 can be set by passing the keyword argument *color* to kwds_quiver_0. The color of the arrows in qu, however, are determined by C_split_0.

The default drawing order is set to: 1. quiver_0 (zorder=1) 2. quiver (zorder=2) 3. scatter (zorder=3) This order can be changed by setting the zorder in kwds_quiver_0, kwds_quiver and/or kwds_scatter. See also http://matplotlib.org/examples/pylab_examples/zorder_demo.html

See also:

```
append_binning_labels_v(), plot_map(), plot_2d(), plot_2d_generator(),
plot_3d()
```

deepgraph.deepgraph.DeepGraph.plot_hist

Plot a histogram (or pdf) of x.

Compute and plot the histogram (or probability density) of x. Keyword arguments are passed to plt.plot. See parameters and np.histogram for details.

Parameters

- **x** (*array_like*) The data from which a frequency distribution is plot.
- **bins** (*int or array_like, optional (default=10)*) If bins is an int, it determines the number of bins to create. If log_bins is True, this number determines the (approximate) number of bins to create for each magnitude. For linear bins, it is the number of bins for the whole range of values. If bins is a sequence, it defines the bin edges, including the rightmost edge, allowing for non-uniform bin widths.
- **log_bins** (bool, optional (default=False)) Whether to use logarithmically or linearly spaced bins.
- **density** (bool, optional (default=False)) If False, the result will contain the number of samples in each bin. If True, the result is the value of the probability *density* function at the bin, normalized such that the *integral* over the range is 1. Note that the sum of the histogram values will not be equal to 1 unless bins of unity width are chosen; it is not a probability *mass* function.
- **floor** (*bool*, *optional* (*default=False*)) Whether to floor the bin edges to the closest integers. Only has an effect if bins is an int.
- **ax** (matplotlib axes object, optional (default=None)) An axes instance to use.

Returns

- **ax** (*matplotlib axes object*) A matplotlib axes instance.
- hist (np.ndarray) The values of the histogram. See density.
- **bin_edges** (*np.ndarray*) The edges of the bins.

deepgraph.deepgraph.DeepGraph.plot_logfile

static DeepGraph.plot_logfile(logfile)

Plot a logfile.

Plot a benchmark logfile created by create_edges or create_edges_ft.

Parameters logfile (*str*) – The filename of the logfile.

Returns obj – Depending on the logfile, return a dict of matplotlib objects with a subset of the following keys: ['fig', 'ax', 'pc_n', 'pc_e', 'cb_n', 'cb_e']

Return type dict

Other Methods

DeepGraph.append_binning_labels_v(col,	Append a column with binning labels of the values in
col_name)	v[col].
DeepGraph.append_cp([directed, connection,	Append a component membership column to v.
])	
<pre>DeepGraph.filter_by_values_v(col, values)</pre>	Keep only nodes in v with features of type col in
	values.
<pre>DeepGraph.filter_by_values_e(col, values)</pre>	Keep only edges in e with relations of type col in
	values.
DeepGraph.filter_by_interval_v(col,	Keep only nodes in v with features of type col in
interval)	interval.
DeepGraph.filter_by_interval_e(col, inter-	Keep only edges in e with relations of type col in
val)	interval.
DeepGraph.update_edges()	After removing nodes in v, update e.

deepgraph.deepgraph.DeepGraph.append_binning_labels_v

DeepGraph.append_binning_labels_v(col, col_name, bins=10, log_bins=False, floor=False, re-

 $turn_bin_edges=False)$ Append a column with binning labels of the values in v[col].

Append a column col_name to v with the indices of the bins to which each value in v [col] belongs to.

If bins is an int, it determines the number of bins to create. If log_bins is True, this number determines the (approximate) number of bins to create for each magnitude. For linear bins, it is the number of bins for the whole range of values. If floor is set True, the bin edges are floored to the closest integer. If return_bin_edges is set True, the created bin edges are returned.

If bins is a sequence, it defines the bin edges, including the rightmost edge, allowing for non-uniform bin widths.

See np.digitize for details.

Parameters

- **col** (*int* or *str*) A column name of v, whose corresponding values are binned and labelled.
- **col_name** (*str*) The column name for the created labels.
- **bins** (*int* or *array_lke*, *optional* (*default=10*)) If bins is an int, it determines the number of bins to create. If log_bins is True, this number determines the (approximate) number of bins to create for each magnitude. For linear bins, it is the number of bins for the whole range of values. If bins is a sequence, it defines the bin edges, including the rightmost edge, allowing for non-uniform bin widths.
- **log_bins** (bool, optional (default=False)) Whether to use logarithmically or linearly spaced bins.
- **floor** (*bool*, *optional* (*default=False*)) Whether to floor the bin edges to the closest integers.

• **return_bin_edges**(bool, optional (default=False)) – Whether to return the bin edges.

Returns

- **v** (*pd.DataFrame*) Appends an extra column col_name to v with the binning labels.
- bin_edges (np.ndarray) Optionally, return the created bin edges.

Examples

First, we need a node table:

```
>>> import pandas as pd
>>> import deepgraph as dg
>>> v = pd.DataFrame({'time': [1,2,12,105,899]})
>>> g = dg.DeepGraph(v)
```

Binning time values with default arguments:

```
>>> bin_edges = g.append_binning_labels_v('time', 'time_l',
... return_bin_edges=True)
```

>>> q.v time time_1 0 1 1 1 2 1 2 12 1 3 105 2 4 899 10

Binning time values with logarithmically spaced bins:

```
>>> bin_edges = g.append_binning_labels_v('time', 'time_l', bins=5,
... log_bins=True,
... return_bin_edges=True)
```

48..

79.,

>>:	> g.v	
	time	time_l
0	1	1
1	2	2
2	12	6
3	105	10
4	899	15

Binning time values with logarithmically spaced bins (floored):

```
>>> bin_edges = g.append_binning_labels_v('time', 'time_l', bins=5,
... log_bins=True, floor=True,
... return_bin_edges=True)
```

>>> g.v

```
time
          time_1
      1
                 1
1
       2
                 2
2
     12
                 5
3
                9
    105
4
    899
               14
```

deepgraph.deepgraph.DeepGraph.append_cp

```
DeepGraph.append_cp(directed=False, connection='weak', col_name='cp', label_by_size=True, con-solidate_singles=False)
```

Append a component membership column to v.

Append a column to v indicating the component membership of each node. Requires scipy.

Parameters

- **directed** (*bool*, *optional* (*default=False*)) If True, then operate on a directed graph: only move from point i to point j along paths csgraph[i, j]. If False, then find the shortest path on an undirected graph: the algorithm can progress from point i to j along csgraph[i, j] or csgraph[j, i].
- connection (*str*, *optional* (*default='weak'*)) One of {'weak','strong'}. For directed graphs, the type of connection to use. Nodes i and j are strongly connected if a path exists both from i to j and from j to i. Nodes i and j are weakly connected if only one of these paths exists. Only has an effect if directed is True
- **col_name** (*str*, *optional* (*default='cp'*)) The name of the appended column of component labels.
- **label_by_size** (*bool*, *optional* (*default=True*)) Whether to rename component membership labels to reflect component sizes. If True, the smallest component corresponds to the largest label, and the largest component corresponds to the label 0 (or 1 if consolidate_singles is True). If False, pass on labels given by scipy's connected_components method directly (faster and uses less memory).
- **consolidate_singles** (bool, optional (default=False)) If True, all singular components (components comprised of one node only) are consolidated un-

der the label 0. Also, all other labels are renamed to reflect component sizes, see label_by_size.

Returns \mathbf{v} – appends an extra column to \mathbf{v} indicating component membership.

Return type pd.DataFrame

deepgraph.deepgraph.DeepGraph.filter_by_values_v

DeepGraph.filter_by_values_v(col, values)

Keep only nodes in v with features of type col in values.

Remove all nodes from v (and their corresponding edges in e) with feature(s) of type col not in the list of features given by values.

Parameters

- **col** (*str* or *int*) A column name of v, indicating the type of feature used in the filtering.
- **values** (*object* or *array_like*) The value(s) indicating which nodes to keep.

Returns

- v (*pd.DataFrame*) update ∨
- e (pd.DataFrame) update e

deepgraph.deepgraph.DeepGraph.filter_by_values_e

DeepGraph.filter_by_values_e(col, values)

Keep only edges in e with relations of type col in values.

Remove all edges from e with relation(s) of type col not in the list of relations given by values.

Parameters

- **col** (*str or int*) A column name of e, indicating the type of relation used in the filtering.
- values (object or array_like) The value(s) indicating which edges to keep.

Returns e – update e

Return type pd.DataFrame

deepgraph.deepgraph.DeepGraph.filter_by_interval_v

DeepGraph.filter_by_interval_v(col, interval, endpoint=True)

Keep only nodes in v with features of type col in interval.

Remove all nodes from v (and their corresponding edges in e) with features of type col outside the interval given by a tuple of values. The endpoint is included, if endpoint is not set to False.

Parameters

• **col** (*str or int*) – A column name of v, indicating the type of feature used in the filtering.

- **interval** (*tuple*) A tuple of two values, (value, larger_value). All nodes outside the interval are removed.
- endpoint (bool, optional (default=True)) False excludes the endpoint.

Returns

- **v** (*pd.DataFrame*) update v
- **e** (*pd.DataFrame*) update e

deepgraph.deepgraph.DeepGraph.filter_by_interval_e

DeepGraph.filter_by_interval_e (col, interval, endpoint=True)

Keep only edges in e with relations of type col in interval.

Remove all edges from e with relations of type col outside the interval given by a tuple of values. The endpoint is included, if endpoint is not set to False.

Parameters

- **col** (*str* or *int*) A column name of e, indicating the type of relation used in the filtering.
- **interval** (*tuple*) A tuple of two values, (value, larger_value). All edges outside the interval are removed.
- endpoint (bool, optional (default=True)) False excludes the endpoint.

Returns e – update e

Return type pd.DataFrame

deepgraph.deepgraph.DeepGraph.update_edges

DeepGraph.update_edges()

After removing nodes in v, update e.

If you deleted rows from v, you can remove all edges associated with the deleted nodes in e by calling this method.

Returns e – update e

Return type pd.DataFrame

1.4.2 The Functions Module

deepgraph.functions

Connector Functions

```
great_circle_dist
cp_node_intersection
cp_intersection_strength
hypergeometric_p_value
```

Selector Functions

1.5 Contact

1.5.1 Email

Please feel free to contact me if you have questions or suggestions regarding DeepGraph:

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Dominik Traxl <dominik.traxl@posteo.org>
```

1.5.2 Authors

Deepgraph was written as part of a PhD thesis in physics by Dominik Traxl at Humboldt University Berlin, the Berstein Center for Computational Neuroscience and the Potsdam Institute for Climate Impact Research.

CHAPTER 2

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