"Install a pip package in the current Jupyter kernel\n", "Memo to self: Install modules with 'sys.executable' and then restart kernel"

```
In [1]: import sys
         #!{sys.executable} -m pip install --upgrade statsmodels
         #****** Start importing python modules
In [17]:
         import time
         import os # For
         import pandas as pd
         import numpy as np
         from functools import partial # For partial functions
         from scipy.interpolate import CubicSpline
         import statsmodels.api as sm
         import scipy.cluster.hierarchy as spc
         from matplotlib import pyplot as plt
         from math import sqrt
         #****** Start GUI modules
         import io
         import traitlets
         import ipywidgets as widgets
         from IPython.display import display
         from tkinter import Tk, filedialog
         #***** End GUI modules
         #****** End importing python modules
         # Start
```

Start importing data First ask the user to select the json file containing

```
In [3]: btn_upload = widgets.FileUpload(
          accept = '*.json',
          multiple = False
)
display(btn_upload)
```

FileUpload(value=(), accept='*.json', description='Upload')

Extract the content of the json-file and convert the json to a data frame

```
In [4]: stringContent = btn_upload.value[0]['content'].tobytes().decode("utf-8")
    df = pd.read_json(io.StringIO(stringContent),orient = 'index')
    print(f'top part is {df.head()}')
    print(f'bottom part is {df.tail()}')
```

```
top part is
                                series_0 series_1 series_2 series_3 series_4
                                            0.0 0.009272 -0.943774
2022-06-01 00:00:00
                         0.0 -0.012866
2022-06-01 00:01:00
                         0.1 0.106740
                                            0.1 -0.004306 -0.734072
2022-06-01 00:02:00
                         0.2 0.209939
                                            0.2 0.005588 3.272961
2022-06-01 00:03:00
                         0.3 0.293082
                                            0.3 -0.005641 0.832609
2022-06-01 00:04:00
                         0.4 0.391814
                                             0.4 -0.006649 -2.225293
                    series 5 series 6
2022-06-01 00:00:00
                         NaN
                                   NaN
2022-06-01 00:01:00
                         NaN
                                   NaN
2022-06-01 00:02:00
                         NaN
                                   NaN
2022-06-01 00:03:00
                       228.0
                                   1.0
2022-06-01 00:04:00
                                   1.0
                         NaN
bottom part is
                                   series_0
                                               series 1
                                                          series_2 series_3 s
eries 4
2022-06-02 23:55:00
                       287.5 287.498961 287.323143 -0.165455 2.017107
2022-06-02 23:56:00
                       287.6 287.599466 287.600000 -0.010588 -1.490229
2022-06-02 23:57:00
                       287.7 287.719646 287.700000 -0.005317 -0.667750
2022-06-02 23:58:00
                       287.8 287.790522 287.800000 0.005404 -3.185175
2022-06-02 23:59:00
                       287.9 287.910746 287.900000 -0.006259 -0.760504
                    series_5 series_6
2022-06-02 23:55:00
                      2375.0
                                   3.0
2022-06-02 23:56:00
                         NaN
                                   3.0
2022-06-02 23:57:00
                         NaN
                                   3.0
2022-06-02 23:58:00
                         NaN
                                   3.0
2022-06-02 23:59:00
                         NaN
                                   3.0
```

End importing dataWe start by imputing missing values using cubic splines

```
In [18]: # v needs to be a single column/vector
         def find index missing(v,include nonmissing = True):
             index_missing_values = [index for index in range(len(v)) if np.isnan(v[index
             returnValue = index_missing_values
             if(include nonmissing):
                 index nonmissing = [index for index in range(len(v)) if index not in ind
                 returnValue = index missing values, index nonmissing
             return returnValue
         # By default 'extrapolate' is set to False since the behavior of cubic splines o
         # and last spline can sometimes be erratic
         def imputeCubicSpline(x,y,extrapolate = False):
             index_missing_values,index_nonmissing = find_index_missing(
                 y,include_nonmissing=True)
             if len(index_missing_values) > 0 :
                 cs = CubicSpline(x[index nonmissing],y[index nonmissing],extrapolate = e
                 y = cs(x)
             return y
         # If 'timeCol' is None, then the index is assumed to be a timestamp. Oth
         # This function imputes missing values by fitting a cubic spline to the non-miss
         # If "imputationCols" is a string (single column), then it is converted to a list
         # I 'imputationCols" is None, then every column in the dataframe is imputed
         def df_impute_cubic(df,imputationCols=None,timeCols = None,crop = False,**kwargs
             if type(imputationCols) is str:
                 imputationCols = [imputationCols]
             elif imputationCols is None:
                 imputationCols = list(df.columns)
```

```
#Datetimes need to be converted to unix time
listOfTimes = None
if type(timeCols) is str:
    listOfTimes = df[timeCol].tolist()
elif timeCols is None:
    listOfTimes = list(df.index)

#
timeValues = np.asarray([ts.timestamp() for ts in listOfTimes])

#
imputer = partial(imputeCubicSpline,timeValues,**kwargs)
df[imputationCols] = df[imputationCols].apply(imputer)

#
if crop:
    df = df.dropna()
    return df
```

Do the imputation

```
In [19]: df = df_impute_cubic(df,crop=True)
```

For single numerical columns x and y (of equal length) we choose aboluste value of the correlation og x and y as our distance metrix i,e,

$$dist(x, y) = 1 - |corr(x, y)|.$$

(1)

When either x or y have missing values we choose to calculate the correlation as the

The reason why we are using the absolute value of the ccrrelation is h For gorups of colums we choose method = 'complete' as our distance metric From documentation: method='complete' assigns

$$d(u,v) = \max(dist(u[i],v[j])).$$

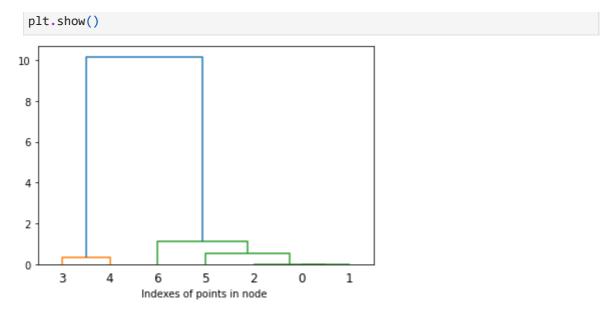
(2)

for all points (in our case columns) `i` in cluster u and `j` in cluster `v`., This is also known by the Farthest Point Algorithm or Voor Hees Algorithm.Description algorithm: 1. Calculate (1) defined above for each pair of columns in the data frame ("condensed values") 2. Do Agglomerative Hiearchical Clustering based We don't want to calculate the correlation of the same two vectors more than once. To avoid that we first calculate (1) defined above for each pair of columns in the data frame ("condensed" and the used that to and pass the "condensed values" (each distance

```
In [47]:
    def CCorDistance(u,v,max_lag = None):
        # Use default value suggested in documentation of sm.tsa.stattools.acf
        centered_u = u - np.mean(u) # Demean
        centered_v = v - np.mean(v) # Demean
        cc_full = np.correlate(a=centered_u, v=centered_v,mode = 'full')
        cc_full = cc_full/ (len(u) * np.std(u) * np.std(v)) # Normalization
        # Only negative lags are used
        cc_non_positive_lags = cc_full[range(len(v)),]
        squared_cc = [cc_non_positive_lags[ind,]**2 for ind in range(cc_non_positive #
        D = sqrt((1.0-squared_cc[-1]) /sum(squared_cc[:-1]))
    return D

#
def condensedSet(df):
    listDists = []
```

```
# col1 ie every column except for the last one
             # col2 ie every column except for the first
             for i in range(df.shape[1]-1):
                 for j in range(i+1,df.shape[1]):
                     listDists.append(CCorDistance(df.iloc[:,i].to_numpy(),df.iloc[:,j].t
             listDists = np.asarray(listDists)
             return listDists
         def clusterByCrossCorr(df,max_intra_dist = None):
             print('Er nå inne i clusterByCrossCorr')
             pdist condensed = condensedSet(df)
             # if max_intra_dist is not provided then st the max distance within the same
             if max_intra_dist is None:
                 max_intra_dist = np.quantile(pdist_condensed,0.2)
             #
             print(f'type(pdist condensed) er {type(pdist condensed)} og pdist condensed.
             print(f'max_intra_dist er {max_intra_dist}')
             print('pdist condensed er')
             print(pdist_condensed)
             linkage = spc.linkage(pdist_condensed, method='complete')
             idx = spc.fcluster(linkage, max_intra_dist, 'distance')
             # Put the cluster information into a data frame. The first column is the nam
             # The second columns is an indicator that says which cluster group each colu
             groupingFrame = pd.DataFrame(list(df.columns),columns=['Series'])
             groupingFrame['Group'] = list(idx)
             clusterDict = {}
             for group in list(set(list(idx))):
                 clusterDict['_'.join(['group',str(group)])] = [
                     list(df.columns)[col] for col in list(groupingFrame[groupingFrame['6
                     1
             return linkage,clusterDict
In [48]: linkage,clusterGroups = clusterByCrossCorr(df)
        Er nå inne i clusterByCrossCorr
       type(pdist_condensed) er <class 'numpy.ndarray'> og pdist_condensed.shape er (2
       1,)
       max_intra_dist er 0.5522145670160576
       pdist condensed er
        [5.51152745e-06 3.43892418e-04 1.01662004e+01 2.64099058e+00
        5.52214567e-01 1.12485465e+00 3.43907633e-04 1.01657290e+01
        2.64099706e+00 5.52215090e-01 1.12486036e+00 9.36371247e+00
        2.64029720e+00 5.52227870e-01 1.12490892e+00 3.45256371e-01
        1.74282065e+00 1.68013211e+00 1.46002755e+00 1.36709347e+00
        6.57015666e-01]
In [49]: print(f' The cluster groups outputted from the algorithm are \n{clusterGroups}')
        The cluster groups outputted from the algorithm are
        {'group_1': ['series_3', 'series_4'], 'group_2': ['series_0', 'series_1', 'series
       _2'], 'group_3': ['series_5'], 'group_4': ['series_6']}
In [50]: spc.dendrogram(linkage)
         plt.xlabel("Indexes of points in node ")
```



Thee first grouping is of "series_0','series_1' and 'series_2'. The distances between these 3 are so small that it is not visible in the dendrogram above.