# Week 11 Lab (Hierarchical Clustering)

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#### **About The Data**

We'll be using the Credit Card Dataset from kaggle for this lab, but feel free to follow along with your own dataset. We will develop a customer segmentation to define marketing strategy. The sample Dataset summarizes the usage behavior of about 9000 active credit card holders during the last 6 months. The file is at a customer level with 18 behavioral variables:

- CUSTID: Identification of Credit Card holder (Categorical) • BALANCE : Balance amount left in their account to make purchases (
- BALANCEFREQUENCY: How frequently the Balance is updated, score between 0 and 1 (1 = frequently updated, 0 = not
- frequently updated) • PURCHASES: Amount of purchases made from account
- ONEOFFPURCHASES: Maximum purchase amount done in one-go
- INSTALLMENTSPURCHASES: Amount of purchase done in installment
- CASHADVANCE: Cash in advance given by the user PURCHASESFREQUENCY: How frequently the Purchases are being made, score between 0 and 1 (1 = frequently purchased,
- 0 = not frequently purchased) ONEOFFPURCHASESFREQUENCY: How frequently Purchases are happening in one-go (1 = frequently purchased, 0 = not frequently purchased)
- PURCHASESINSTALLMENTSFREQUENCY: How frequently purchases in installments are being done (1 = frequently done, 0 = not frequently done)
- CASHADVANCEFREQUENCY: How frequently the cash in advance being paid CASHADVANCETRX: Number of Transactions made with "Cash in Advanced"
- PURCHASESTRX: Numbe of purchase transactions made
- CREDITLIMIT: Limit of Credit Card for user • PAYMENTS : Amount of Payment done by user
- MINIMUM\_PAYMENTS: Minimum amount of payments made by user
- PRCFULLPAYMENT : Percent of full payment paid by user
- TENURE: Tenure of credit card service for user
- **About Hierarchical clustering** Hierarchical clustering is a method which seeks to build a hierarchy of clusters. Strategies for hierarchical clustering generally fall
- into two types:

moves down the hierarchy.

• Agglomerative: This is a "bottom-up" approach: each observation starts in its own cluster, and pairs of clusters are merged as one moves up the hierarchy. • Divisive: This is a "top-down" approach: all observations start in one cluster, and splits are performed recursively as one

- In general, the merges and splits are determined in a greedy manner, and the results of hierarchical clustering are usually presented in a dendrogram.
- **Agglomerative Clustering**

 Compute the proximity matrix Let each data point be a cluster

### Repeat: Merge two closest clusters and update the proximity matrix until 1/ K cluster remains

The algorithm goes as follows:

- For example, say we have six data points {a,b,c,d,e,f}
  - In the initial step, we consider all the six data points as individual clusters as shown in the image below.
- The first step is to determine which elements to merge in a cluster. Usually, we want to take the two closest elements, according to the chosen distance. We construct a distance matrix at this stage, where the number in the i-th row j-th column is the distance

between the i-th and j-th elements. Then, as clustering progresses, rows and columns are merged as the clusters are merged and the distances updated.

Computation of Proximity/Distance Matrix

following: • The maximum distance between elements of each cluster (also called complete-linkage clustering) The minimum distance between elements of each cluster (also called single-linkage clustering)

To calculate the proximity between two clusters, we need to define the distance between them. Usually the distance is one of the

Note: Euclidean, Manhattan, Mahalanobis, etc. distance formulas can be used when calculating distances for each of the above.

The mean distance between elements of each cluster (also called average linkage clustering)

- **Implementation**
- Similarly to the K Means lab, we'll skip the data exploration portion and jump to implementation, but you're welcome to explore

import matplotlib.pyplot as plt

rcParams['figure.figsize'] = 15, 5

0.818182

0.909091

The sum of all intra-cluster variance.

this data, or your own if working with a different dataset.

import numpy as np import pandas as pd

cc df.head()

**1** 3202.467416

0

Out[5]: BALANCE

40.900749

BALANCE FREQUENCY

ONEOFF PURCHASES

CASH ADVANCE TRX

PRC FULL PAYMENT

dtype: int64

TENURE

INSTALLMENTS PURCHASES

CASH ADVANCE FREQUENCY

ONEOFF PURCHASES FREQUENCY

PURCHASES INSTALLMENTS FREQUENCY

PURCHASES FREQUENCY

PURCHASES

CASH ADVANCE

import seaborn as sns

sns.set\_style('darkgrid')

from matplotlib import rcParams

Let's first load the data into a pandas DataFrame. MINIMUM\_PAYMENTS and CREDIT\_LIMIT are missing a few values, so we'll go

BALANCE BALANCE\_FREQUENCY PURCHASES ONEOFF\_PURCHASES INSTALLMENTS\_PURCHASES CASH\_ADVANCE PURCH

0.00

0.00

773.17

0.000000

6442.945483

95.4

0.0

ahead and fill them with their respective column mean. We'll also go ahead and drop CUST\_ID while we're at it since it serve's us no valuable information. cc\_df = pd.read\_csv('cc.csv') cc df.drop('CUST ID', axis = 1, inplace=True) cc df['MINIMUM PAYMENTS'].fillna(value=cc df['MINIMUM PAYMENTS'].mean(), inplace = True) cc\_df['CREDIT\_LIMIT'].fillna(value=cc\_df['CREDIT\_LIMIT'].mean(), inplace = True)

2 2495.148862 0.000000 1.000000 773.17 0.0 1499.00 **3** 1666.670542 0.636364 1499.00 0.0 205.788017 817.714335 1.000000 16.00 0.0 0.000000 16.00 Checking to see that there are no more missing values: cc df.isnull().sum()

95.40

0.00

0 0

0 0

0

0

0

0

0

0 0

PURCHASES TRX 0 0 CREDIT LIMIT PAYMENTS 0 MINIMUM PAYMENTS 0

We'll now standardize and normalize the data as follows: from sklearn.preprocessing import StandardScaler, normalize scaler = StandardScaler() scaled\_df = scaler.fit\_transform(cc\_df) normalized df = normalize(scaled df) # Converting the scaled data back into a pandas DataFrame normalized df = pd.DataFrame(data=normalized df) Next (optional, but recommended), we will reduce the dimensions of the data using PCA from sklearn.decomposition import PCA pca = PCA(n\_components = 2) X principal = pca.fit transform(normalized df)

X\_principal.head()

**0** -0.489826 -0.679678

-0.482374 -0.092114

**4** -0.563289 -0.481914

40

30

20

10

X principal = pd.DataFrame(X\_principal) X principal.columns = ['P1', 'P2']

**P2** 

-0.518791 0.545010 0.330885 0.268980

import scipy.cluster.hierarchy as shc plt.title('visualising the data') Dendrogram = shc.dendrogram((shc.linkage(X principal, method ='ward'))) 60 50

Note: The cell below might take a while to run.

from sklearn.cluster import AgglomerativeClustering from sklearn.metrics import silhouette score # where we'll save scores for later plotting silhouette\_scores = [] # testing different cluster values in range [2,8) for n cluster in range(2, 8): silhouette\_scores.append(silhouette\_score(X\_principal, AgglomerativeClustering(n\_clusters = n\_cluster).fit\_predict(X\_

# Creating bar graph to compare the results. You can use a line plot if you prefer (similar to K Means lab)

Number of clusters

Next, we'll use a dendrogram to visualize the linkage of the reduced data (X\_principal) using **method='ward'**.

visualising the data

Silhouette Score 0.2 0.1

plt.bar(x=range(2, 8), height=silhouette scores)

plt.xlabel('Number of clusters') plt.ylabel('Silhouette Score')

# creating and fitting model

Out[13]: AgglomerativeClustering(n\_clusters=3)

# Visualizing the clustering

agg.fit(X\_principal)

agg = AgglomerativeClustering(n clusters=3)

plt.show()

0.4

0.3

0.0

0.50

In [14]:

We can also determine the optimal number of clusters using silhouette score:

plt.scatter(X\_principal['P1'], X\_principal['P2'], c = AgglomerativeClustering(n clusters = 3).fit predict(X principal)) plt.show()

We'll go ahead and build and visualize a clustering model for n\_clusters = 3

1.00 0.75

0.50

0.75

1.00

0.25 0.00 -0.25

-0.50-0.75

Congrats! (2) Now try repeating the lab steps on your own data for practice.