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E-tivity 3: Clustering and Manifold Learning

Use this notebook to complete Tasks 1 and 2 in E-tivity3.

Import Python Modules

```
In [1]: import pandas as pd
import numpy as np
import matplotlib.pyplot as plt

from sklearn import manifold
from sklearn import cluster
from sklearn import preprocessing
```

- In [2]: # Using a global random seed for reproducible results when doing multiple runs
 np.random.seed(42)
- In [3]: # Load the dataset into pandas dataframe
 df = pd.read_csv("loans_dataset_et3.csv")
- In [4]: # Look at the first 5 data samples
 df.head()
- Out[4]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_Histo
0	2483	2466.0	90	180	0
1	4917	0.0	130	360	0
2	4106	0.0	40	180	1
3	3859	3300.0	142	180	1
4	6417	0.0	157	180	1

In [5]: # Descriptive statistics of the data
df.describe()

Out[5]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_
count	332.000000	332.000000	332.000000	332.000000	332.000
mean	5201.093373	1495.508795	140.882530	341.710843	0.97891
std	4584.815491	1982.742932	75.544237	61.651497	0.14388
min	645.000000	0.000000	17.000000	60.000000	0.00000
25%	2912.750000	0.000000	100.000000	360.000000	1.00000
50%	3858.500000	1211.500000	128.000000	360.000000	1.00000
75%	5818.250000	2250.000000	162.000000	360.000000	1.00000
max	39999.000000	20000.000000	600.000000	480.000000	1.00000
4					

Observation:

There are no missing values in the dataset. The data contains 332 examples and 5 features/attributes

The dataset features have different scales with a ApplicantIncome and CoapplicantIncome being of larger order of magnitude compared to LoanAmount and Loan_Amount_Term. Credit_History seems to be a categorical attribute with binary outcome.

Decision: Scaling is necessary for this dataset before applying clustering algorithm

In [6]: df.shape
Out[6]: (332, 5)

Task 1 (CS5062)

Cluster loans_dataset_et3.csv with the k-Means clustering algorithm and visualise the clustering. Use an appropriate feature-scaling technique and pick a value of k suggested by MDS and t-SNE visualisations of the dataset.

Experiment with various values of the parameters of k-Means. **Aim at producing a meaningful clustering** that allows you to describe the clusters with words. Write a short conclusion about the characteristics of the clusters (max 250 words).

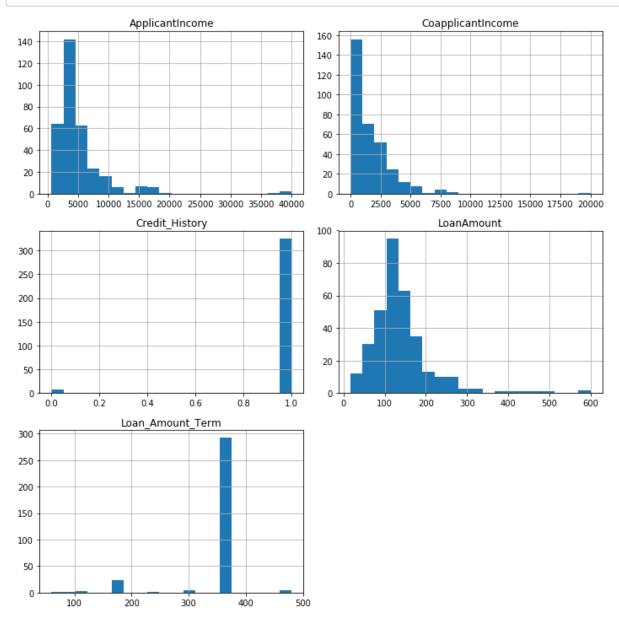
Tests:

- 1. Use an appropriate feature-scaling technique
- 2. pick a value of k suggested by MDS and t-SNE visualisations of the dataset.
- 3. Experiment with various values of the parameters of k-Means
- 4. Aim at producing a meaningful clustering that allows you to describe the clusters with words.

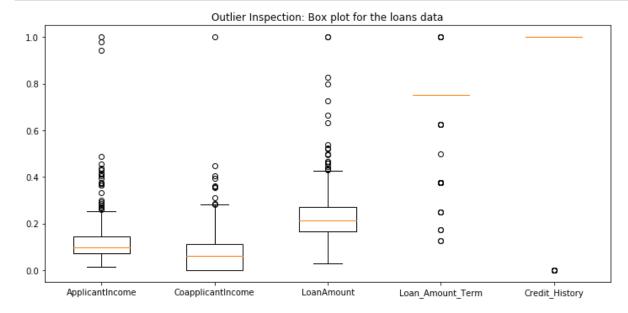
Reference: https://scikit-learn.org/stable/auto_examples/manifold/plot_compare_methods.html (<a href="https://scikit-learn.org/stable/auto_examples/manifold/plot_compare_methods.htm

```
In [7]: X = np.array(df.values, dtype=float)
In [8]: X.shape
Out[8]: (332, 5)
```

In [9]: # Check the distribution of Loan dataset
 df.hist(figsize=(10,10), bins=20)
 plt.tight_layout()
 plt.show()



```
In [10]: # Check if there are any outlier as it may affect the scaling algorithm
    plt.figure(figsize=(10,5))
    plt.boxplot([df[i]/df[i].max() for i in df.keys()], labels=df.keys())
    plt.title('Outlier Inspection: Box plot for the loans data')
    plt.tight_layout()
    plt.show()
```



Observation on the Distribution and Box Plots:

From the boxplot we can see that there are some potential outliers in the dataset and this can affect the scaling outcome and Proper scaling technique needs to be used to counter the outliers so that we can get a good clustering outcome.

Scaling the data:

Scaling should be applied when there are input features on a different order of magnitude, or even large differences in the same order of magnitude, than other features. We need to take into account outliers and their effects on scaling. If the data contains many outliers(as in this case), scaling using the mean and variance of the data is likely to not work very well.

Performed a study on the various approaches to do scaling in scikit-learn:

- StandardScaler: Removes the mean and scales the data to unit variance
- MinMaxScaler: Rescales the data set such that all feature values are in the range [0, 1]
- MaxAbsScaler: The absolute values are mapped in the range [0, 1], same as MinMaxScaler if we have only +ve data
- RobustScaler: The centering and scaling statistics of this scaler are based on percentiles and are therefore not influenced by a few number of very large marginal outliers
- PowerTransformer: applies a power transformation to each feature to make the data more Gaussianlike
- QuantileTransformer: (Gaussian: saturates outliers & uniform: same as robust scaler) it smooths out
 unusual distributions and is less influenced by outliers than scaling methods.

Decision on scaling: I have decide to use the QuantileTransformer with uniform scaling

Note: StandardScaler, MinMaxScaler and MeanAbsScaler is very sensitive to the presence of outliers and hence should not be used here.

References:

- https://scikit-learn.org/stable/auto_examples/preprocessing/plot_all_scaling.html (https://scikit-learn.org/stable/auto_examples/preprocessing/plot_all_scaling.html)
- https://scikit-learn.org/stable/modules/preprocessing.html (https://scikit-learn.org/stable/modules/preprocessing.html)

```
In [11]: # Scale the data with appropriate scaler which can handle ouliers
    quantile_scaler = preprocessing.QuantileTransformer(output_distribution='unifo
    rm')
    quantile_uniform_data = quantile_scaler.fit_transform(X)
```

In [12]: # Reusing the plotting method from Lab 3 notebook # Create an array of colors to distinguish the clusters during plotting colors = np.array(['orange', 'blue', 'lime', 'khaki', 'pink', 'green', 'purpl e', 'maroon', 'aqua', 'black', 'yellow', 'gray', 'magenta', 'cy an']) # points - a 2D array of (x,y) coordinates of data points # labels - an array of numeric labels in the interval [0..k-1], one for each p oint # centers - a 2D array of (x, y) coordinates of cluster centers # title - title of the plot def clustering_scatterplot(points, labels, centers, title): # plot the examples, i.e. the data points n_clusters = np.unique(labels).size for i in range(n clusters): h = plt.scatter(points[labels==i,0], points[labels==i,1], c=colors[i%colors.size], label = 'cluster '+str(i)) # plot the centers of the clusters if centers is not None: plt.scatter(centers[:,0], centers[:,1], c='r', marker='*', s=400) plt.title(title) plt.legend(loc='best') plt.xlabel('x') plt.ylabel('y')

Manifold techniques used for K-Means:

MDS:

The purpose of the MDS is to find a low-dimensional representation of the data (here 2D) in which the distances respect well the distances in the original high-dimensional space, unlike other manifold-learning algorithms, it does not seeks an isotropic representation of the data in the low-dimensional space.

Reference: https://scikit-learn.org/stable/modules/manifold.html (https://scikit-learn.org/stable/modules/modules/modules/modules/modules/modules/modules/modules/modules/modules/modules/module

T-SNE (T-distributed Stochastic Neighbor Embedding):

T-SNE tries to map high-dimensional into low-dimensional objects by preserving the neighbourhood relationship structure in spite of a trade-off consisting of a misclassification around the far objects. Drawback of T-SNE is that its computationally very expensive and 10 times slower than MDS

T-SNE Parameters:

- n_components: default (equal to 2), it is the number of the ending dimensions.
- Perplexity: It is the main parameter. The higher perplexity, the more the the number of neighbours. A
 good rule is setting it up between 5-50.
 - If Perplexity is too small we will not capture enough neighbours; on the contrary, an excessive number will approach values too far.
- · learning rate: A range between 10 and 1000 is usually recommended

References:

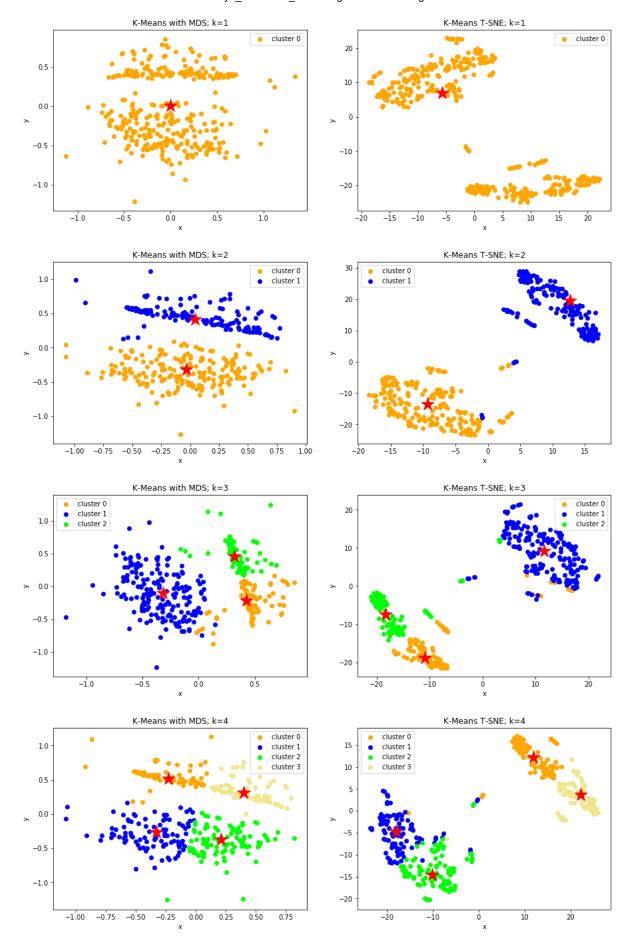
- 1. https://www.deeplearningitalia.com/stochastic-neighbor-embedding-sne-its-correction-in-t-sne/ (https://www.deeplearningitalia.com/stochastic-neighbor-embedding-sne-its-correction-in-t-sne/)
- 2. https://scikit-learn.org/stable/auto_examples/manifold/plot_compare_methods.html (https://scikit-learn.org/stable/auto_examples/manifold/plot_compare_methods.html)

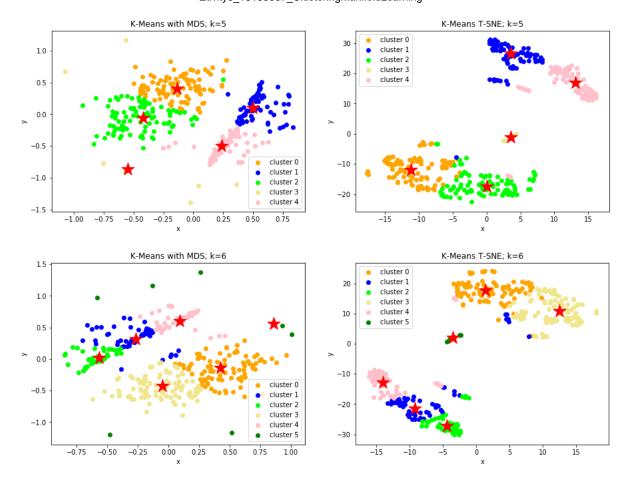
```
In [13]:
          def kmeans clustering(data, k):
             # Apply K-Means Clustering technique
             kmeans = cluster.KMeans(n clusters=k, n init=10, max iter=500).fit(data)
             # append the cluster centers to the dataset
             data_and_centers = np.r_[data, kmeans.cluster_centers_]
             # Apply the MDS transformation
             XY_mds_coordinates = manifold.MDS(n_components=2).fit_transform(data_and_c
         enters)
             plt.subplots(1, 2, figsize=(15, 5))
             plt.subplot(1, 2, 1)
             # Plot the MDS clusters
             clustering scatterplot(points=XY mds coordinates[:-k,:],
                                     labels=kmeans.labels ,
                                     centers=XY_mds_coordinates[-k:,:],
                                     title='K-Means with MDS; k={}'.format(k))
             # Apply the t-SNE transformation to project both the data and the k-Means
          cluster centers to a 2D space
             XY_tsne_coordinates = manifold.TSNE(n_components=2,
                                                  perplexity=30,
                                                  learning rate=200).fit transform(data
         and centers)
             plt.subplot(1, 2, 2)
             # Plot the t-SNE clusters
             clustering_scatterplot(points=XY_tsne_coordinates[:-k,:],
                                     labels=kmeans.labels ,
                                     centers=XY tsne coordinates[-k:,:],
                                     title='K-Means T-SNE; k={}'.format(k))
             return kmeans, data_and_centers
```

```
In [14]: # Select the maximum and minimium number of clusters for experiment
    max_num_clusters = 7
    min_num_clusters = 1
```

```
In [15]: output = dict()

# Call the K-means clustering function
for n in range(min_num_clusters, max_num_clusters):
    output[n] = kmeans_clustering(quantile_uniform_data, n)
    plt.show()
```





Cluster Analysis

The visual inspection of the plots suggest that the most probable number of clusters is 4

In [16]: # Create a copy of dataframe
 df_kmeans = df.copy()

Add the cluster labels of k=4 as an extra column in the original loan datafr
 ame
 df_kmeans['cluster'] = pd.Series(output[4][0].labels_, index=df.index)

In [17]: df_kmeans.head()

Out[17]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_Histo
0	2483	2466.0	90	180	0
1	4917	0.0	130	360	0
2	4106	0.0	40	180	1
3	3859	3300.0	142	180	1
4	6417	0.0	157	180	1
4					

In [19]: # Compute the mean value of each attribute for each cluster
df_kmeans.groupby('cluster').mean()

Out[19]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_
cluster					
0	9338.506667	16.720000	198.493333	336.800000	0.9600
1	5096.419355	3085.333333	176.150538	343.225806	0.9784
2	2640.200000	2181.967579	103.694737	347.242105	0.9789
3	4370.855072	14.956522	81.927536	337.391304	1.0000
4					

Observation from the clusters using K-means:

Cluster number	description
Cluster 0	Customers with very high Applicant Income group with a very high loan amount with low Coapplicant Income
Cluster 1	Customers with average Applicant Income group and very high Co-applicant Income with above average loan amount
Cluster 2	Customers with very low Applicant Income and above average Co-applicant Income
Cluster 3	Customers with below average Applicant Income and very low Co-applicant Income with very low Loan Amount and a Credit History of 1

Task 2 (CS5062)

Plot the sum of squared distances from the data points (in loans_dataset_et3.csv) to the centers of the k-Means clusters for various values of k. Use the Elbow method to pick the best value of k. Attempt also another method for determining the best value of k. If these methods suggest a different value of k (from the one you used in Task 1), then compute a new k-Means clustering.

Use another clustering algorithm (from the ones available in scikit-learn) for the same dataset and an appropriate manifold-learning technique to visualise it.

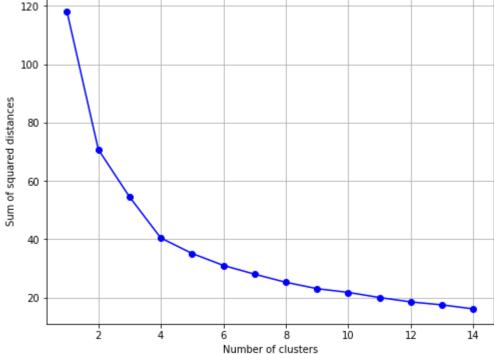
Discuss the differences between the k-Means clustering and the clustering found by the second algorithm you have tried in words (max 250 words) and by visualising the clustering's with an appropriate manifold-learning technique.

Plotting the Sum of Squared distances and applying Elbow method

```
In [20]: SSE = []
         # Create a range of clusters
         num clusters = range(1,15)
         # For loop to run K-means for each cluster vakue 'k' and find the SSE
         for k in num_clusters:
             km = cluster.KMeans(n clusters=k, n init=10, max iter=500).fit(quantile un
         iform data)
             SSE.append(km.inertia_)
             print("SSE for {} cluster = {}".format(k, km.inertia_))
         plt.figure(figsize=(8,6))
         plt.plot(num_clusters, SSE, 'bo-')
         plt.xlabel('Number of clusters')
         plt.ylabel('Sum of squared distances')
         plt.title("K-Means Clustering: Elbow Criterion Plot", fontsize=14, fontweight=
         'bold')
         plt.grid()
         plt.show()
```

```
SSE for 1 cluster = 117.93565366310898
SSE for 2 cluster = 70.78355526826287
SSE for 3 cluster = 54.512640531014995
SSE for 4 cluster = 40.44126607418319
SSE for 5 cluster = 35.137093923613115
SSE for 6 cluster = 31.04218934834786
SSE for 7 cluster = 28.02647710231204
SSE for 8 cluster = 25.244839279946724
SSE for 9 cluster = 23.057725126307222
SSE for 10 cluster = 21.765492180429746
SSE for 11 cluster = 20.019258184016927
SSE for 12 cluster = 18.502588173669555
SSE for 13 cluster = 17.49460196168556
SSE for 14 cluster = 16.11280021037343
```

K-Means Clustering: Elbow Criterion Plot



Observation on Elbow plot:

Usually an 'elbow' is apparent where the reduction in sum of square errors(SSE) levels off as the number of clusters increses. This indicates that increasing the number of clusters does not improve the clustering.

The elbow can be seen at the optimal number of clusters, k = 4 where the SSE is not reducing drastically(slope is less) and even if we add more clusters the SSE decay is very slow and not adding any significance to the clustering process

Attempt another method for determining the best value of k and an appropriate manifold-learning technique to visualise it.

Reference for clustering algorithms: https://scikit-learn.org/stable/modules/clustering.html (https://scikit-learn.org/stable/module

Clustering Algorithm Selected: Affinity Propagation

Hyper-parameters:

- preference
 - can be used to control how many data points are selected as exemplars
- damping
 - damping factor dampfact should be at least 0.5 and less than 1
 - Higher damping factors will lead to slower convergence

Manifold techniques used for Affinity Propagation:

- 1. Isomap
- 2. T-SNE

References:

- https://scikit-learn.org/stable/modules/generated/sklearn.cluster.AffinityPropagation.html (https://scikit-learn.org/stable/modules/generated/sklearn.cluster.AffinityPropagation.html (https://scikit-learn.org/stable/modules/generated/sklearn.cluster.AffinityPropagation.html (https://scikit-learn.org/stable/modules/generated/sklearn.cluster.AffinityPropagation.html (https://scikit-learn.org/stable/modules/generated/sklearn.cluster.AffinityPropagation.html)
- https://www.psi.toronto.edu/affinitypropagation/faq.html)
 https://www.psi.toronto.edu/affinitypropagation/faq.html)
- An attempt to use other techniques besides MDS and t-SNE. (https://www.youtube.com/watch?v=j8080l9Pvic))

```
In [21]: def affinity propagation(data, preference):
             # Function to do the clustering using Affinity Propagation
             # Affinity Propagation on various 'preference' values
             af clustering = cluster.AffinityPropagation(damping=0.9,
                                                          max iter=500,
                                                          convergence iter=50,
                                                          preference=preference,
                                                          verbose=True).fit(data)
             af cluster centers = af clustering.cluster centers
             # get the number of clusters
             af clusters = len(af clustering.cluster centers indices )
             # Append the cluster centers to the dataset
             af_data_and_centers = np.r_[data, af_cluster_centers]
             plt.subplots(1, 2, figsize=(15, 5))
             plt.subplot(1, 2, 1)
             # Trying Isomap manifold technique
             XYcoordinates = manifold.Isomap(n neighbors=5, n components=2).fit transfo
         rm(af_data_and_centers)
             # plot the transformed examples and the centers
             clustering scatterplot(points=XYcoordinates[:-af clusters,:],
                                     labels=af_clustering.labels_,
                                     centers=XYcoordinates[-af clusters:,:],
                                     title='Isomap for Affinity Propagation; preference
         {}; clusters={}'.format(
                                         preference, af clusters))
             plt.subplot(1, 2, 2)
             # Apply the t-SNE transformation to project both the data and the k-Means
          cluster centers to a 2D space
             XY tsne coordinates = manifold.TSNE(n components=2,
                                                  perplexity=30,
                                                  learning_rate=200).fit_transform(af_da
         ta_and_centers)
             # Plot the t-SNE clusters
             clustering_scatterplot(points=XY_tsne_coordinates[:-af_clusters,:],
                                     labels=af clustering.labels ,
                                     centers=XY tsne coordinates[-af clusters:,:],
                                     title='T-SNE for Affinity Propagation; preference
         {}; clusters={}'.format(
                                         preference, af clusters))
             plt.show()
             return af_clustering, af_clusters
```

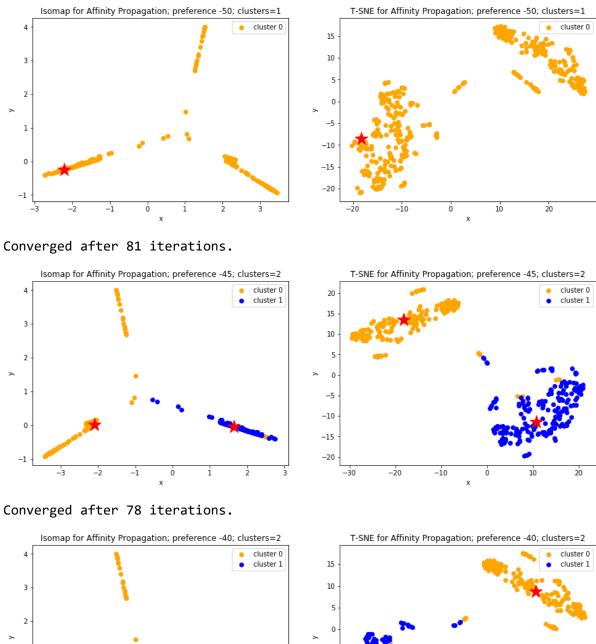
```
In [22]: # Store the scaled data in X_scaled
    X_scaled = quantile_uniform_data

# Hyperparameter choice for preference
    max_preference = 0
    min_preference = -50

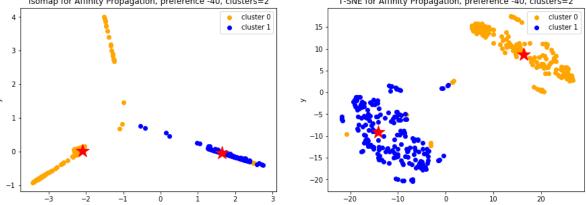
# Jump 5 steps per preference value
    preference_range = range(min_preference, max_preference, 5)

result = dict()
    num_clusters=[]
    for pref in preference_range:
        result[pref], af_cluster = affinity_propagation(X_scaled, pref)
        num_clusters.append(af_cluster)
```

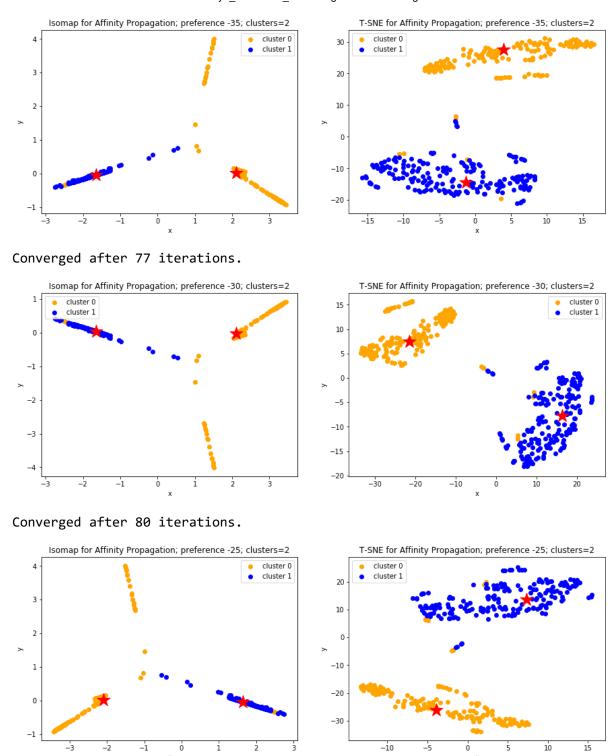
Converged after 76 iterations.



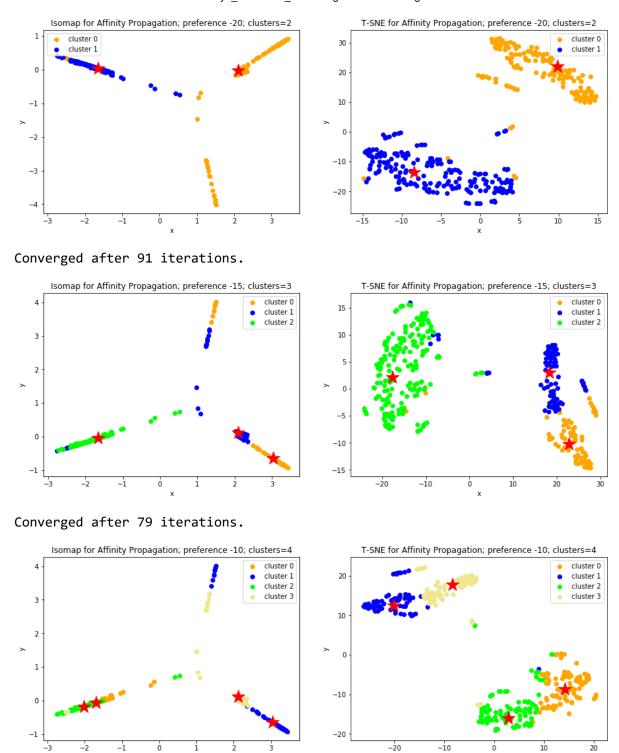




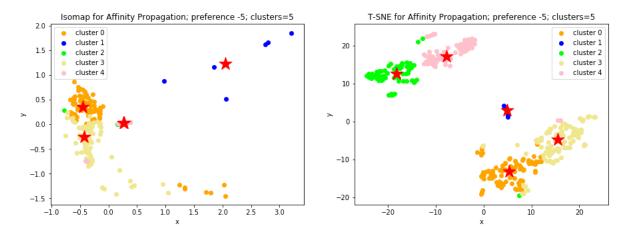
Converged after 76 iterations.



Converged after 82 iterations.



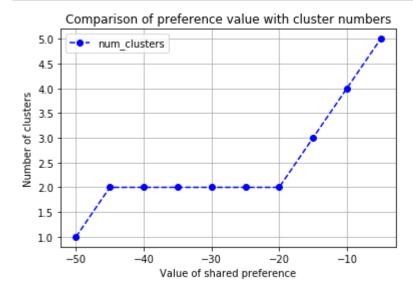
Converged after 107 iterations.



In [23]: # Store the preference and cluster number from Affinity propagation
 clusters_df = pd.DataFrame({'preference_range': preference_range, 'num_cluste
 rs':num_clusters})
 clusters_df

Out[23]:

	preference_range	num_clusters
0	-50	1
1	-45	2
2	-40	2
3	-35	2
4	-30	2
5	-25	2
6	-20	2
7	-15	3
8	-10	4
9	-5	5



Observations from Affinity Propagation clustering technique:

The optimal number of clusters after using the Affinity Propagation clustering algorithm using various preference values is seen to be 2. The visual inspection using T-SNE manifold technique however shows that 4 clusters are more appropriate.

The value of preference plays a very important role in picking up the number of clusters in the Affinity Propagation clustering technique. For preference values -20 to -45 we can see that the number of clusters found are always 2. For other values of preference the number of clusters are not stable and change very quickly. Even with 0 preference value we will have a lot of clusters (almost all data points are clusters in that case).

General observation was:

- low preferences leading to small numbers of clusters
- · high preferences leading to large numbers of clusters

preference value	num_clusters
-50	1
-45 to -20	2
-15	3
-10	4
-5	5

Cluster Analysis: Affinity Propagation clustering technique

In [25]: # Create a copy of dataframe
df_afprop = df.copy()

Using perplexity of -10 in Affninity propagation results which gives 4 clust ers

df_afprop['cluster'] = pd.Series(result[-10].labels_, index=df.index)

df_afprop.head()

Out[25]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_Histo
0	2483	2466.0	90	180	0
1	4917	0.0	130	360	0
2	4106	0.0	40	180	1
3	3859	3300.0	142	180	1
4	6417	0.0	157	180	1

In [26]: df_afprop.mean()

Out[26]: ApplicantIncome 5201.093373 CoapplicantIncome 1495.508795 LoanAmount 140.882530

dtype: float64

Out[27]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit_
cluster					
0	9338.506667	16.720000	198.493333	336.800000	0.9600
1	5096.419355	3085.333333	176.150538	343.225806	0.9784
2	2640.200000	2181.967579	103.694737	347.242105	0.9789
3	4370.855072	14.956522	81.927536	337.391304	1.0000

In [28]: # Affinity Propagation clusters
df_afprop.groupby('cluster').mean()

Out[28]:

	ApplicantIncome	CoapplicantIncome	LoanAmount	Loan_Amount_Term	Credit
cluster					
0	2583.336957	2205.825217	103.315217	346.826087	0.9782
1	4454.112676	14.535211	82.957746	338.028169	1.0000
2	5007.389474	3054.326316	173.715789	343.578947	0.9789
3	9420.972973	32.162162	201.013514	336.486486	0.9594
4			1		•

Observation from the clusters using Affinity Propagation:

The results of Affinity propagation clusters are very similar to the K-Means clustering results for 4 clusters.

Cluster number	Cluster Description
Cluster 0	Customers with very low Applicant Income and above average Co-applicant Income
Cluster 1	Customers with below average Applicant Income and very low Co-applicant Income with very low Loan Amount and a Credit History of 1
Cluster 2	Customers with average Applicant Income group and very high Co-applicant Income with above average loan amount
Cluster 3	Customers with very high Applicant Income group with a very high loan amount with low Coapplicant Income

Differences between the k-Means clustering and the clustering found by Affinity Propagation algorithm

K-means algorithms clusters the data into n clusters of equal variance minimizing the within-cluster Euclidean distance.

- · The algorithm requires the samples and the number of clusters as input
- · k-means assumes that clusters are convex shaped and forms cluster of even size

Affinity Propagation algorithm forms clusters by sending messages between pairs of instances until convergence

- · The algorithm choses the clusters based on the data
- Affinity propagation finds "exemplars", members of the input set that are representative of clusters
 where damping damps the responsibility and availability messages, preference controls the number of
 exemplars
- · It forms many clusters of uneven size

Silhouette score analysis to understand the optimal number of clusters

The silhouette analysis measures how well an observation is clustered and it estimates the average distance between clusters.

The silhouette_score gives the average value for all the samples. Silhouette coefficients range between -1 and 1, with 1 indicating dense, well separated clusters. This gives a perspective into the density and separation of the formed clusters

Reference:

- 1. https://stackoverflow.com/questions/19197715/scikit-learn-k-means-elbow-criterion)

 (https://stackoverflow.com/questions/19197715/scikit-learn-k-means-elbow-criterion)
- https://chrisalbon.com/machine_learning/clustering/evaluating_clustering/ (https://chrisalbon.com/machine_learning/clustering/evaluating_clustering/)
- 3. https://www.datanovia.com/en/lessons/cluster-validation-statistics-must-know-methods/#silhouette-coefficient)

 methods/#silhouette-coefficient)

```
In [29]: from sklearn.metrics import silhouette_score

# Loop to find the silhouette score for k=2 to k=10 number of clusters
for k in range(2, 11):

# Apply K-Means clustering on the scaled data
kmeans = cluster.KMeans(n_clusters=k, n_init=10, max_iter=500).fit(X_scale
d)

# Get the labels
label = kmeans.labels_

# Calculate the silhoutte score
sil_coeff = silhouette_score(X_scaled, label, metric='euclidean')

print("For n_clusters={}, The Silhouette Coefficient is {}".format(k, sil_coeff))
```

```
For n_clusters=2, The Silhouette Coefficient is 0.3958331519475405
For n_clusters=3, The Silhouette Coefficient is 0.3717353085258766
For n_clusters=4, The Silhouette Coefficient is 0.3685989117677958
For n_clusters=5, The Silhouette Coefficient is 0.3828607398313568
For n_clusters=6, The Silhouette Coefficient is 0.34162019857901854
For n_clusters=7, The Silhouette Coefficient is 0.36408926084722965
For n_clusters=8, The Silhouette Coefficient is 0.33267424065672535
For n_clusters=9, The Silhouette Coefficient is 0.3411891279954565
For n_clusters=10, The Silhouette Coefficient is 0.35690598569914117
```

The maximum score for silhouette is for cluster number 2 and cluster number 5

Lets visualize the Silhouette

The silhouette plot displays a measure of how close each point in one cluster is to points in the neighboring clusters.

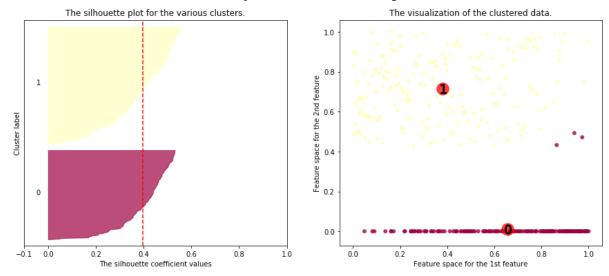
Reference: http://www.awesomestats.in/python-cluster-validation/ (htt

```
In [30]: from sklearn.metrics import silhouette samples
         cluster range = range(2, 7)
         # Loop to find the silhouette score for k=2 to k=6 number of clusters
         for n_clusters in cluster_range:
             # Create a subplot with 1 row and 2 columns
             fig, (ax1, ax2) = plt.subplots(1, 2)
             fig.set_size_inches(15, 6)
             # The (n clusters+1)*10 is for inserting blank space between silhouette pl
         ots of individual clusters,
             # to demarcate them clearly.
             ax1.set ylim([0, len(X scaled) + (n clusters + 1) * 10])
             # Initialize the clusterer with n clusters value
             clusterer = cluster.KMeans(n_clusters=n_clusters, n_init=10, max_iter=500)
             cluster_labels = clusterer.fit_predict(X_scaled)
             # Calulate the average silhouette score
             silhouette_avg = silhouette_score(X_scaled, cluster_labels)
             print("For n clusters =", n clusters, "The average silhouette score is :",
         silhouette avg)
             # Compute the silhouette scores for each sample
             sample silhouette values = silhouette samples(X scaled, cluster labels)
             y lower = 10
             for i in range(n clusters):
                 # Aggregate the silhouette scores for samples belonging to cluster i,
          and sort them
                 ith cluster silhouette values = sample silhouette values[cluster label
         s == i]
                 ith_cluster_silhouette_values.sort()
                 size_cluster_i = ith_cluster_silhouette_values.shape[0]
                 y_upper = y_lower + size_cluster_i
                 color = plt.cm.Spectral(float(i) / n clusters)
                 ax1.fill_betweenx(np.arange(y_lower, y_upper), 0, ith_cluster_silhouet
         te_values,
                                   facecolor=color, edgecolor=color, alpha=0.7)
                 # Label the silhouette plots with their cluster numbers at the middle
                 ax1.text(-0.05, y_lower + 0.5 * size_cluster_i, str(i))
                 # Compute the new y_lower for next plot
                 y lower = y upper + 10
             ax1.set title("The silhouette plot for the various clusters.")
             ax1.set xlabel("The silhouette coefficient values")
             ax1.set_ylabel("Cluster label")
             # The vertical line for average silhoutte score of all the values
             ax1.axvline(x=silhouette avg, color="red", linestyle="--")
```

```
# Clear the yaxis labels / ticks
   ax1.set_yticks([])
   ax1.set_xticks([-0.1, 0, 0.2, 0.4, 0.6, 0.8, 1])
   # 2nd Plot showing the actual clusters formed
   colors = plt.cm.Spectral(cluster labels.astype(float) / n clusters)
   ax2.scatter(X_scaled[:, 0], X_scaled[:, 1], marker='o', s=30, lw=0.35, alp
ha=0.8, c=colors)
   # Labeling the clusters
   centers = clusterer.cluster_centers_
   # Draw red circles at cluster centers
   ax2.scatter(centers[:, 0], centers[:, 1], marker='o', c="red", alpha=0.75,
s = 300)
   for i, c in enumerate(centers):
        ax2.scatter(c[0], c[1], color='black', marker='$%d$' % i, alpha=1, s=1
75)
   ax2.set title("The visualization of the clustered data.")
   ax2.set xlabel("Feature space for the 1st feature")
   ax2.set_ylabel("Feature space for the 2nd feature")
   plt.suptitle(("Silhouette analysis for KMeans clustering with k={}".format
(n_clusters)),
                 fontsize=15, fontweight='bold')
   plt.show()
```

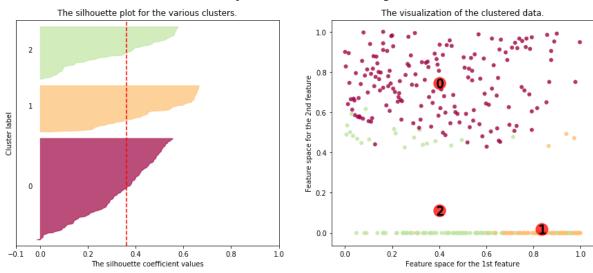
For n_clusters = 2 The average silhouette_score is : 0.3958331519475405

Silhouette analysis for KMeans clustering with k=2



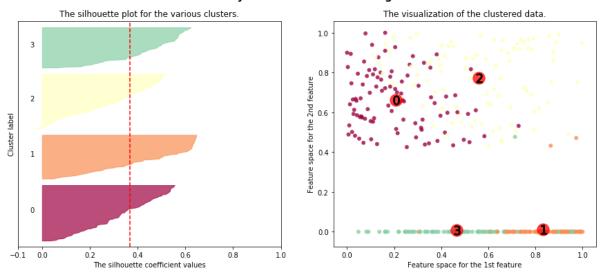
For n_clusters = 3 The average silhouette_score is : 0.36124416254672037

Silhouette analysis for KMeans clustering with k=3



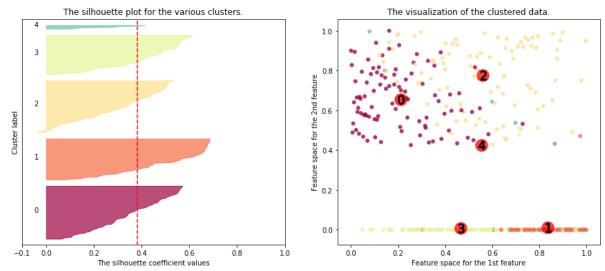
For n_clusters = 4 The average silhouette_score is : 0.3685989117677958

Silhouette analysis for KMeans clustering with k=4



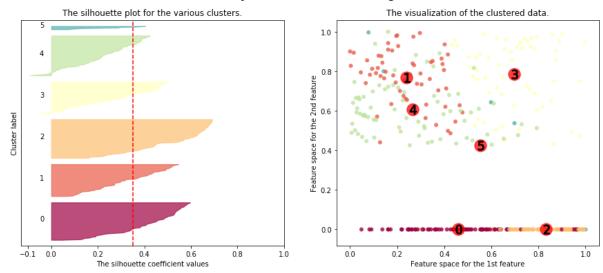
For n_clusters = 5 The average silhouette_score is : 0.3832755211182266

Silhouette analysis for KMeans clustering with k=5



For n_clusters = 6 The average silhouette_score is : 0.3514178223370438

Silhouette analysis for KMeans clustering with k=6



Observation on Silhouette score visualization:

The above plots suggests that the maximum Silhouette score is for K-Means with 2 and 5 clusters. Also there are observations with a negative score (for K-Means with k>2) and hence probably some samples are placed in the wrong cluster for higher number of clusters.

The visual inspection is a bit subjective here but we can see that 2 clusters are more appropriate as the cluster sizes are fairly homogeneous with high Silhouette score and non-negative coefficient values.