ECE 219 Project 2

Tanner Dulay, Loïc Maxwell, Zoë Tucker February 3, 2021

1 Question 1

In this assignment, we explore various methods for clustering of text data. To begin, we use the "20 Newsgroups" dataset, taking the two well-separated classes previously defined ("computer technology" and "recreational activity"). As this project uses unsupervised learning, we do not distinguish between training and test data, and instead use the entire dataset throughout the project.

To start our process, we build the TF-IDF matrix. We use min_df = 3, exclude English stopwords/punctuation, and remove headers and footers, but we do not perform stemming or lemmatization. The resulting TF-IDF matrix is 7882×21909 .

2 Question 2

Next, we apply k-means clustering directly to the TF-IDF data. As we are interested in two classes of documents, we set k=2. The below contingency matrix (Figure 1) shows the results of this.

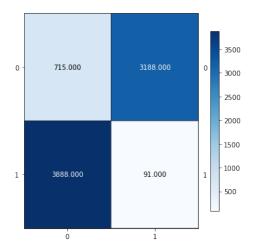


Figure 1: Contingency Matrix for K-Means on TF-IDF Matrix

We then find the metrics we need to evaluate our classification results. We find the following values:

Homogeneity score	0.5600025412974119
Completeness score	0.5716557539454379
V_measure score	0.5657691483417876
Adjusted Rand Index score	0.6327475635535726
Adjusted Mutual Information Score	0.5657289846874574

4 Question 4

As the metrics found in Question 3 do not demonstrate a good clustering result, we now try dimensionality reduction to help our data better fit the assumptions made by k-means algorithms.

We begin by finding the percent of variance the top r principal components can retain for various values of r. Figure 2 shows our findings.

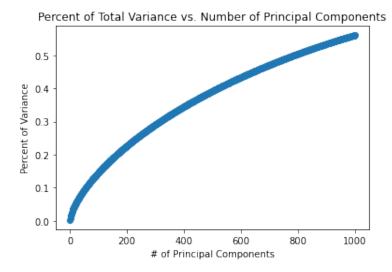


Figure 2: Principal Component Results

Next, we will try to experimentally find the value of r (i.e. the number of dimensions) that we can reduce our data to in order to obtain the best clustering results. Testing $r \in \{1, 2, 3, 5, 10, 20, 50, 100, 300\}$ using both NMF and SVD, we plot the values of the five metrics found in Question 3 for each r. The results are shown in Figures 3 and 4.

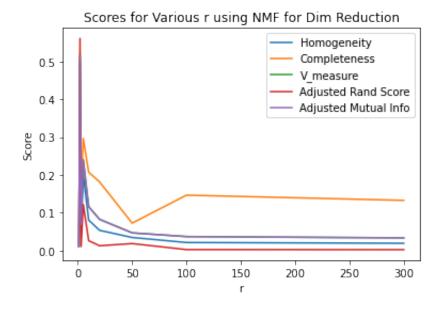


Figure 3: NMF Results

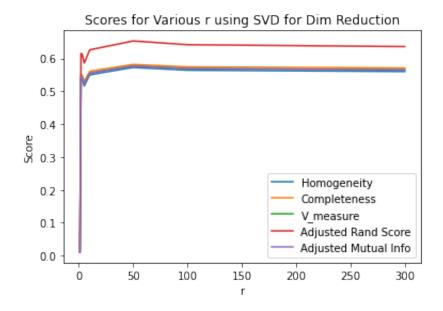


Figure 4: SVD Results

As we can see, the behavior of the measures as r increases is non-monotonic.

This is due to a trade-off between a larger value of r giving a more faithful representation of the data, and a smaller value of r being better suited for k-means clustering. The higher the dimensionality of a dataset, the more difficult it is to meaningfully evaluate the "distance" between data points, which is what k-means relies on. Thus, we expect there to be an optimal value of r that balances these two concerns and maximizes our scores. In our case, we find that r=2 for NMF, and r=50 for SVD.

7 Question 7

Now, we want to visualize our clustering results. We do this by using k-means clustering with our dataset reduced by NMF and SVD with the optimal values of r found in Question 6, then projecting the data into 2-dimensional space with SVD and coloring the points according to their label. Figures 5 and 6 show the results for NMF with r=2. Figures 7 and 8 show the results for SVD with r=50.

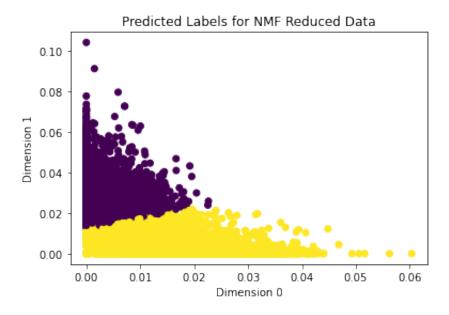


Figure 5: Predicted labels for NMF reduced data

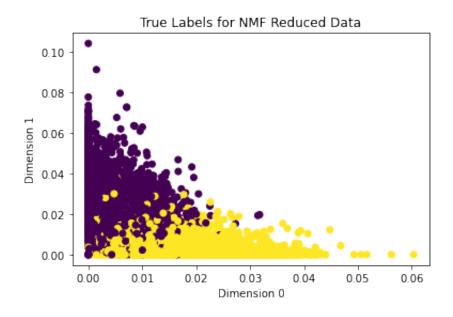


Figure 6: Ground truth labels for NMF reduced data

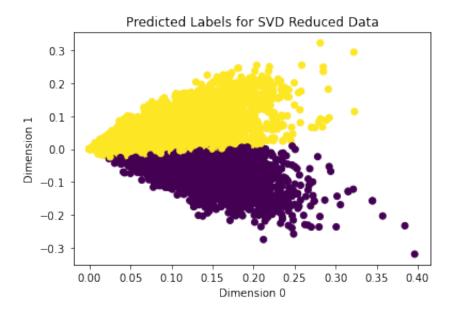


Figure 7: Predicted labels for SVD reduced data

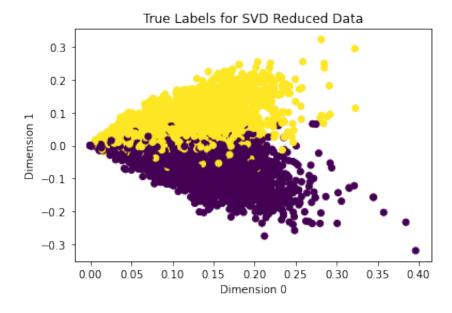


Figure 8: Ground truth labels for SVD reduced data

Because the NMF reduced data was already in 2D, we did not reduce it further with SVD. However, if we were to do SVD reduction, the resulting matrix would still be positive, as it is in Figures 5 and 6. SVD is an orthogonal projection that preserves the angles between feature vectors.

The 2D visualization from Figure 8 shows that the SVD reduced data looks somewhat linearly separable around a value of 0 for Dimension 1, which stems from the second principal component of the TF-IDF matrix. This suggests that this principal component holds important information pertaining to the class that the data belongs to. Although this analysis is restricted to two dimensions, this also suggests that this type of distribution would be good for KMeans clustering. A similar type of separation can be found at the diagonal between dimensions 0 and 1 for the NMF reduced data in Figure 6. This would also make it ideal for KMeans clustering.

However, for both types of reductions, because there is no way to perfectly linearly separate the data there are some better options out there for properly classifying this data. Perhaps a classifier that can learn non-linear behavior in the data, such as a neural network with ReLU/eLU, can perform better on this dataset.

Using SVD with r=50, which gave us the best results on the two classes, we got the following contingency matrix for KMeans on the entire dataset of 20 classes:

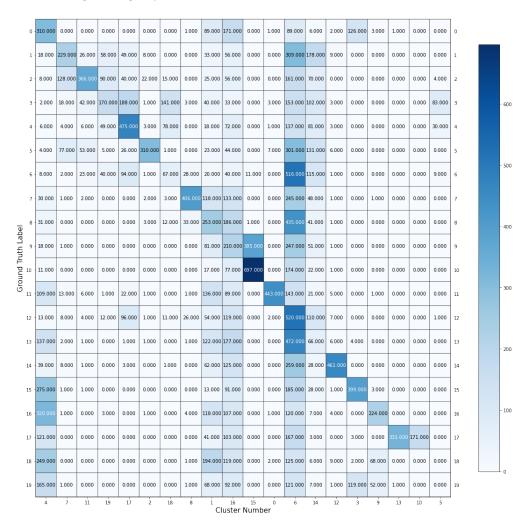


Figure 9: Contingency Matrix for KMeans on 20 Class Data

The respective scores for this clustering are as follows:

Homogeneity score	0.31630458145750995
Completeness score	0.3682552023183464
V_measure score	0.3403086491478565
Adjusted Rand Index score	0.10379799750286293
Adjusted Mutual Information Score	0.3380068278889245

Using the Kullback-Leibler Divergence loss metric for NMF and r=10, we were able to get the following scores:

Homogeneity score	0.37714629232162333
Completeness score	0.3900505160910544
V_measure score	0.3834898799077068
Adjusted Rand Index score	0.18833101982676798
Adjusted Mutual Information Score	0.38146535004967785

There is indeed a noticeable improvement over the performance using ordinary SVD with r=50 (from Q9).

11 Question 11

We compared the performance of KMeans on the dataset reduced with UMAP using various values for n_components. The plots are shown below:

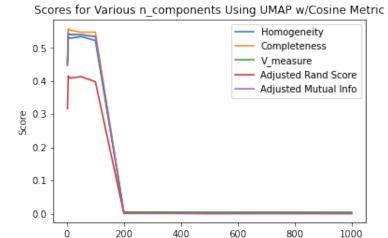


Figure 10: Metric scores for Cosine Metric vs. n_components

n components

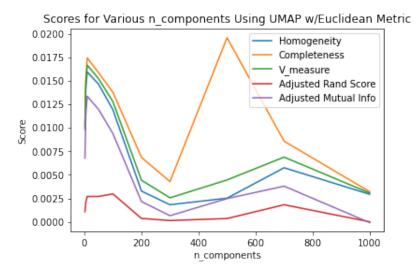


Figure 11: Metric scores for Euclidean Metric vs. ${\tt n_components}$

The best scores using both metrics came from n_{-} components=5:

Cosine Metric	
Homogeneity score	0.5347926404764849
Completeness score	0.5549487009354954
V_measure score	0.5446842655666289
Adjusted Rand Index score	0.4167210932677909
Adjusted Mutual Information Score	0.5431756896373161

Euclidean metric	
Homogeneity score	0.008428056566878686
Completeness score	0.008490790660743977
V_measure score	0.008459307306639803
Adjusted Rand Index score	0.0014762057832842827
Adjusted Mutual Information Score	0.00524781170211466

12 Question 12

Using n_components = 5 as determined above, we get the below contingency matrices – Figure 12 corresponds to the Euclidean metric, and Figure 13 corresponds to the cosine metric.

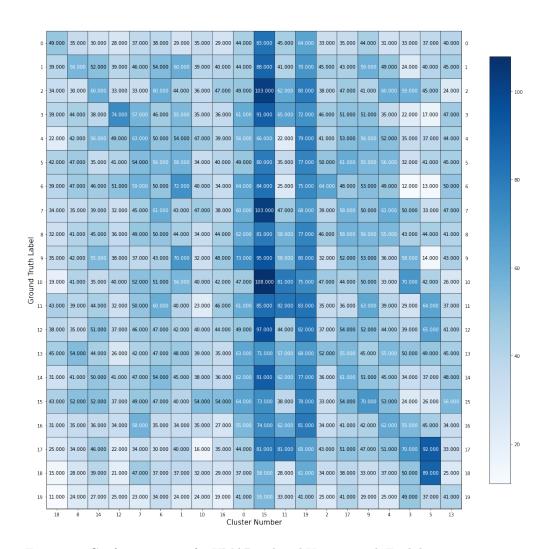


Figure 12: Confusion matrix for UMAP-reduced Kmeans with Euclidean metric

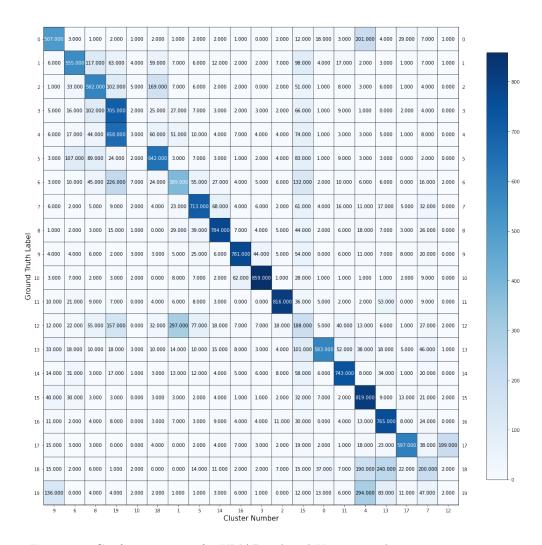


Figure 13: Confusion matrix for UMAP-reduced Kmeans with cosine metric

The cosine matrix is obviously much better, which corresponds with the greatly improved scores using this metric.

Examining the cosine metric more closely, we see that most categories were well identified, but some were more likely to be confused with each other. Tracing the connections, we can see that points with ground truth label 19 were likely to be identified with points labeled 0 or 15, and points from label 12 were misidentified as 1 or 6. In the original dataset, this means that points from talk.religion.misc were identified as alt.atheism or soc.religion.christian, and points from sci.electronics were identified as comp.graphics or misc.forsale. The religion groups form an intuitive grouping, and the confusion between the latter three groups perhaps indicates that selling hardware and electronics was

common on these newsgroups.

13 Question 13

Using the best hyperparameter combination from Q11/12, n_components=5 and the cosine metric, we reduced the Tfidf matrix with UMAP. The scores for agglomerative clustering using 'ward' and 'single' linkage on this data are shown below:

Ward linkage	
Homogeneity score	0.524473300498607
Completeness score	0.5479912976714623
V_measure score	0.53597443687123
Adjusted Rand Index score	0.3940501791969448
Adjusted Mutual Information Score	0.5344316766896055

Single linkage	
Homogeneity score	0.022917457355135236
Completeness score	0.3464156710750703
V_measure score	0.04299081646294147
Adjusted Rand Index score	0.00046761208355549543
Adjusted Mutual Information Score	0.03748616963750802

It's clear that ward linkage performs significantly better on the dataset. Single linkage uses the minimum distance between all points in each cluster to determine whether to merge them or not, so it's a lot more sensitive to outliers and tends to result in uneven cluster sizes. Ward linkage on the other hand, minimizes the variance between pairs of clusters, which can produce more regular cluster sizes and hence better results for our particular dataset.

14 Question 14

In this section we investigate the DBSCAN and HDBSCAN clustering algorithms. We kept UMAP with n_components=5 and the cosine metric as the dimensionality reduction algorithm. The HDBSCAN min_cluster_size was set to 100, and other hyperparameters were tuned to improve the performance of each of these algorithms. We performed hyperparameter tuning on DBSCAN and HDBSCAN for epsilon values of [0.01,0.05,0.1,0.3,0.5,0.7,0.9,1], metric types of [Euclidean, cosine, Manhattan] (if applicable), and minimum sample values of [10, 20, 50, 70, 100]. The following table depicts the scores for each clustering algorithm for the hyperparameter combination that resulted in the best average score across all 5 clustering metrics:

DBSCAN: Epsilon = 0.5, Metric = Manhattan, Minimum samples = 10	
Homogeneity score	0.5407607595568981
Completeness score	0.45471096739764455
V_measure score	0.49401673890037173
Adjusted Rand Index score	0.2967260971944622
Adjusted Mutual Information Score	0.47920988791262803

HDBSCAN: Epsilon = 0.3, Metric = Euclidean, Minimum samples = 10	
Homogeneity score	0.43889528158878366
Completeness score	0.5288819318072416
V_measure score	0.47970500064413385
Adjusted Rand Index score	0.2565568616115164
Adjusted Mutual Information Score	0.4780526456871693

DBSCAN was the better performing clustering algorithm. The associated contingency matrix is shown below:

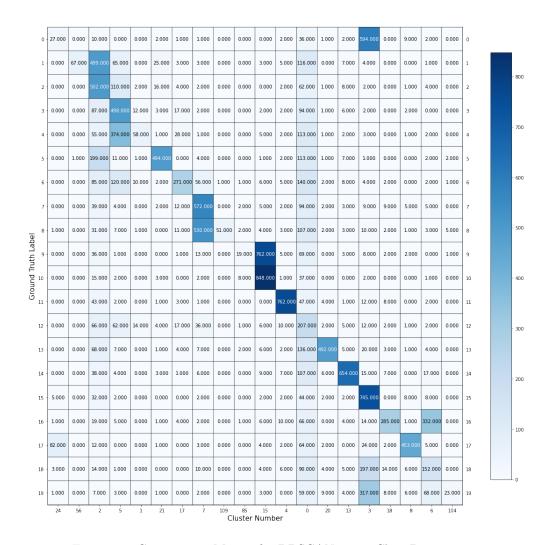


Figure 14: Contingency Matrix for DBSCAN on 20 Class Data

The contingency matrix indicates that there must be at least 96 clusters since that's the largest number shown on the x-axis ticks. By looking at the number of unique labels in the trained DBSCAN model, we are able to determine that there are actually 168 clusters formed by the algorithm (excluding the -1 labels). This number is high because of our low value for the mimimum samples hyperparameter, so most of these cluters contain few data points. However, most of the data points were captured by the larger clusters shown in the contingency matrix, which resulted in acceptable clustering metric scores.

A label of -1 indicates that the data point was classified as noise, and was not assigned to any cluster. This cluster label is represented as cluster number 0 in the contingency matrix. We can tell because this cluster contains many

data points from each of the 20 classes.

16 Question 16

We used the "read_csv" function of the Pandas library to acquire the BBC dataset as a dataframe. After converting the written class labels to numerical target values for each document, we then used our established pipeline to vectorize (excluding stopwords/punctuation) then normalize the data and analyze the dimensions of the resulting TF-IDF matrix. As with the 20 Newsgroups dataset, we then reduced dimensionality with UMAP and searched for the optimal parameter configuration for multiple clustering algorithms. Starting with KMeans, we tested how our five aforementioned accuracy metrics vary with the number of principle components used in clustering for both Euclidean (Figure 15) and Cosine distance metrics (Figure 16). The best combination was n_c components = 10 with the Euclidean distance metric.

Scores for Various n_components Using UMAP w/Euclidean Metric

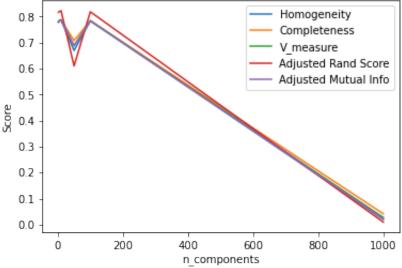


Figure 15: Variation in accuracy metrics by number of components for Kmeans Clustering and Euclidean distance

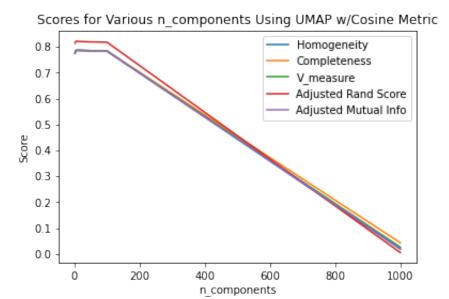


Figure 16: Variation in accuracy metrics by number of components for Kmeans Clustering and Cosine distance

We then tested the accuracy of the DBSCAN clustering algorithm for both distance metrics. We cycled through values from .01 to 1 for epsillon and from 1 to 50 for the minimum number of points per cluster to find the best combination of hyperparameters for our analysis. We used 3-dimensional surface plots to visualize how these parameters in concert affected our five accuracy metrics and the estimated number of clusters generated by DBSCAN (Figure 17).

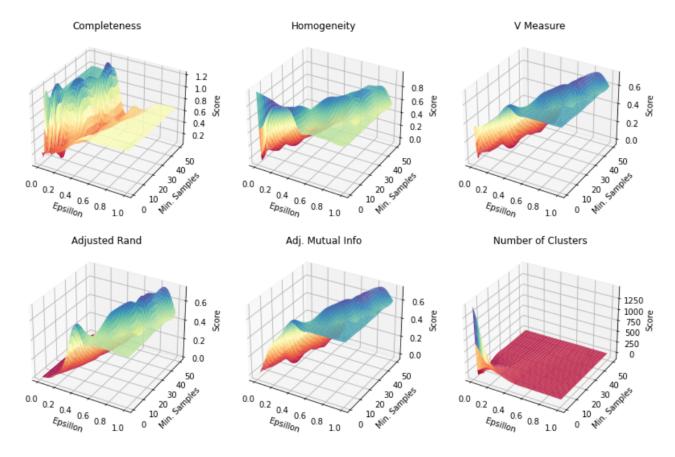


Figure 17: Variation in accuracy metrics corresponding to changes in epsilon and minimum cluster metrics for DBSCAN clustering of UMAP reduced data with Cosine distance metric (Euclidean metric results visually similar)

Avoiding the low values of epsilon and minimum samples that resulted in excessively high numbers of estimated clusters, we found the combination of parameters that maximized the average of the accuracy metrics to be 0.8 for epsilon and 35 for minimum samples. We then compared the results of each of our protocols (Kmeans with Euclidean/Cosine, DBSCAN with Euclidean/Cosine) to find the best reduction and clustering technique for our data. We report the five accuracy metrics for each protocol below:

Accuracy Metrics for DBSCAN, $Metric = Euclidean$	
Homogeneity score	0.7896799513244787
Completeness score	0.667782538810624
V_measure score	0.7236336939202158
Adjusted Rand Index score	0.7108732847743878
Adjusted Mutual Information Score	0.7221232251101021

Accuracy Metrics for DBSCAN, Metric = Cosine	
Homogeneity score	0.7896799513244787
Completeness score	0.667782538810624
V_measure score	0.7236336939202158
Adjusted Rand Index score	0.7108732847743878
Adjusted Mutual Information Score	0.7221232251101021

Accuracy Metrics for KMeans, Metric = Euclidean	
Homogeneity score	0.7874360944663693
Completeness score	0.7874929063651126
V_measure score	0.7874644993910623
Adjusted Rand Index score	0.8229763601775667
Adjusted Mutual Information Score	0.786746935945491

Accuracy Metrics for KMeans, Metric = Cosine	
Homogeneity score	0.7863911146778318
Completeness score	0.7863053396249526
V_measure score	0.7863482248123015
Adjusted Rand Index score	0.8206748429226303
Adjusted Mutual Information Score	0.785626959948192

The best scoring protocol was Kmeans with Euclidean distance metric and n_components set to 10. The contingency matrix for this clustering is shown in Figure 18 below

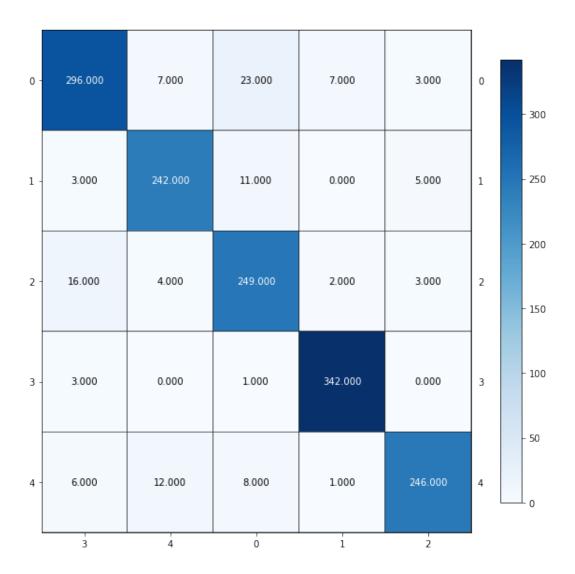


Figure 18: Contingency Matrix for optimal clustering of BBC News data

Our pipeline and parameter selections produce a fairly good clustering through both DBSCAN and KMeans. One factor that may affect the accuracy of our analysis is the relatively small size of the BBC dataset when compared to the 20 Newsgroups dataset. With more data, our clustering would likely become more accurate.

Project_2

February 3, 2021

```
[88]: #01
      import nltk
      from nltk import pos_tag
      from sklearn.feature_extraction.text import TfidfTransformer
      from sklearn.feature_extraction.text import CountVectorizer
      from string import punctuation
      from nltk.corpus import stopwords
      from matplotlib import pyplot as plt
      from sklearn.datasets import fetch_20newsgroups
      categories = ['comp.graphics', 'comp.os.ms-windows.misc',
      'comp.sys.ibm.pc.hardware', 'comp.sys.mac.hardware',
      'rec.autos', 'rec.motorcycles',
      'rec.sport.baseball', 'rec.sport.hockey']
      #Vectorizer
      analyzer = CountVectorizer().build_analyzer()
      tfidf_transformer = TfidfTransformer()
      combined_stopwords = set.union(set(stopwords.words('english')),set(punctuation))
      def stem_rmv_punc_nolem(doc):
          return (word for word in analyzer(doc) if word not in combined_stopwords and_
       →not word.isdigit())
      #Download dataset
      dataset = fetch_20newsgroups(subset='all', categories = categories, shuffle =__
       →True, random_state = 0, remove=('headers', 'footers'))
      count_vect = CountVectorizer(min_df=3,analyzer=stem_rmv_punc_nolem,_
       ⇔stop_words='english')
      X_counts = count_vect.fit_transform(dataset.data)
      X_tfidf = tfidf_transformer.fit_transform(X_counts)
```

```
print("X_tfidf shape: ", X_tfidf.shape)
```

X_tfidf shape: (7882, 21909)

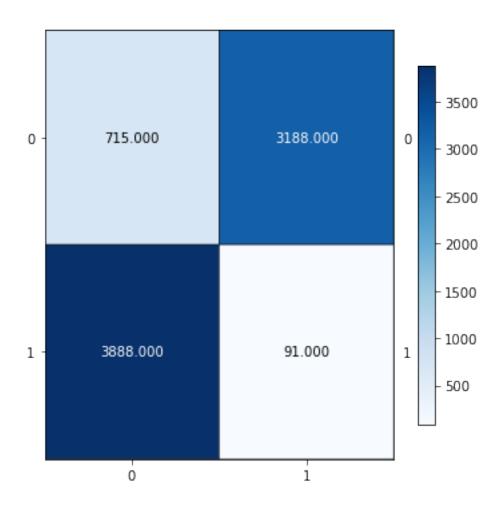
```
[89]: #Q2
#Get labels to verify k means accuracy

from sklearn.cluster import KMeans
from plotmat import plot_mat
from sklearn.metrics.cluster import contingency_matrix

dataset_targets_bin = dataset.target.copy()
for i in range(len(dataset_targets_bin)):
    if dataset.target[i] in [0,1,2,3]:
        dataset_targets_bin[i] = 0
    else:
        dataset_targets_bin[i] = 1

kmeans_clf = KMeans(n_clusters=2, max_iter=1500, random_state=0, n_init=50)
kmeans_clf.fit(X_tfidf)
km_cont_mat = contingency_matrix(dataset_targets_bin, kmeans_clf.labels_)

plot_mat(km_cont_mat, size=(5,5))
```



Homogeneity score: 0.5600025412974119 Completeness score: 0.5716557539454379 V_measure score: 0.5657691483417876

Adjusted Rand Index score: 0.6327475635535726

```
[91]: #Q4

from sklearn.decomposition import TruncatedSVD, NMF
import numpy as np

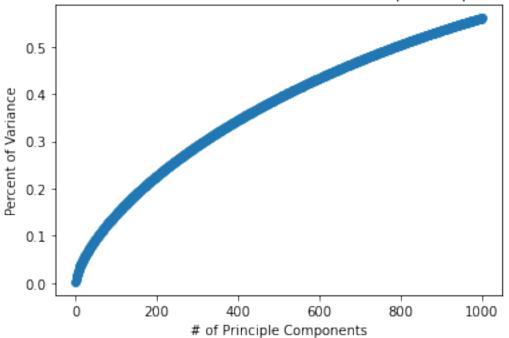
variances = []
r_vals = np.arange(1, 1001)

trunc_svd = TruncatedSVD(n_components=1000, random_state=0)
X_svd = trunc_svd.fit_transform(X_tfidf)

variances = trunc_svd.explained_variance_ratio_

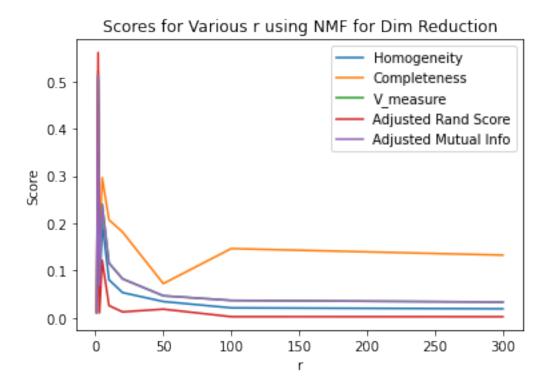
plt.scatter(r_vals, np.cumsum(variances))
plt.xlabel("# of Principle Components")
plt.ylabel("Percent of Variance")
plt.title("Percent of Total Variance vs. Number of Principle Components")
plt.show()
```

Percent of Total Variance vs. Number of Principle Components



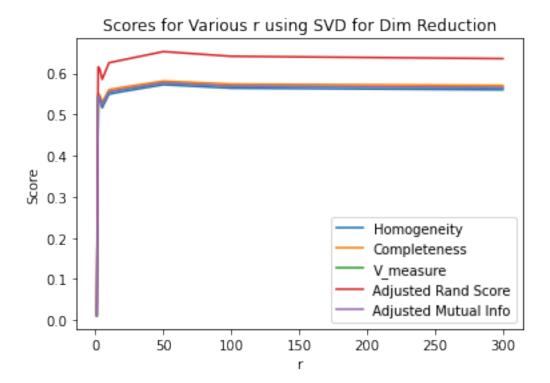
```
[92]: #Q5
r_vals_5 = [1,2,3,5,10,20,50,100,300]
metric_scores_svd = []
```

```
metric_scores_nmf = []
      for r in r_vals_5:
          nmf_model = NMF(n_components=r, init='random', random_state=0, max_iter=1000)
          nmf_red = nmf_model.fit_transform(X_tfidf)
          svd_red = X_svd[:,:r]
          kmeans_clf.fit(nmf_red)
          metric_scores_nmf.append((metrics.homogeneity_score(dataset_targets_bin,_
       ⇔kmeans_clf.labels_),
                                    metrics.completeness_score(dataset_targets_bin,__
       →kmeans_clf.labels_),
                                    metrics.v_measure_score(dataset_targets_bin,__
       →kmeans_clf.labels_),
                                    metrics.adjusted_rand_score(dataset_targets_bin,_
       →kmeans_clf.labels_),
                                    metrics.
       →adjusted_mutual_info_score(dataset_targets_bin, kmeans_clf.labels_)))
          kmeans_clf.fit(svd_red)
          metric_scores_svd.append((metrics.homogeneity_score(dataset_targets_bin,_
       →kmeans_clf.labels_),
                                    metrics.completeness_score(dataset_targets_bin,__
       ⇔kmeans_clf.labels_),
                                    metrics.v_measure_score(dataset_targets_bin,__
       →kmeans_clf.labels_),
                                    metrics.adjusted_rand_score(dataset_targets_bin,__
       →kmeans_clf.labels_),
                                    metrics.
       →adjusted_mutual_info_score(dataset_targets_bin, kmeans_clf.labels_)))
[93]: plt.plot(r_vals_5, metric_scores_nmf)
      plt.legend(("Homogeneity", "Completeness", "V_measure", "Adjusted Rand Score", u
      →"Adjusted Mutual Info"))
      plt.xlabel("r")
      plt.ylabel("Score")
      plt.title("Scores for Various r using NMF for Dim Reduction")
      plt.show()
```



```
[94]: plt.plot(r_vals_5, metric_scores_svd)
plt.legend(("Homogeneity","Completeness","V_measure", "Adjusted Rand Score",

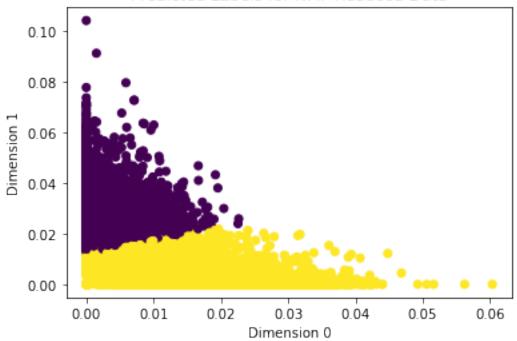
→"Adjusted Mutual Info"))
plt.xlabel("r")
plt.ylabel("Score")
plt.title("Scores for Various r using SVD for Dim Reduction")
plt.show()
```

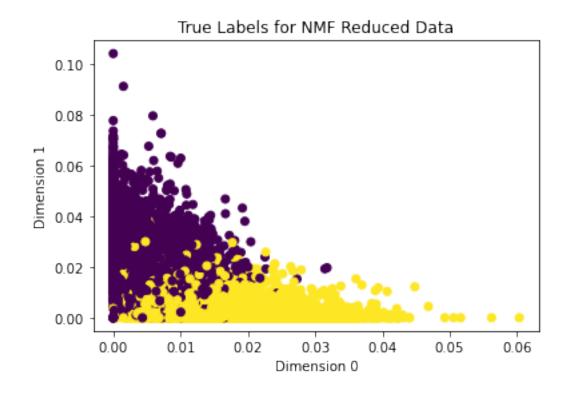


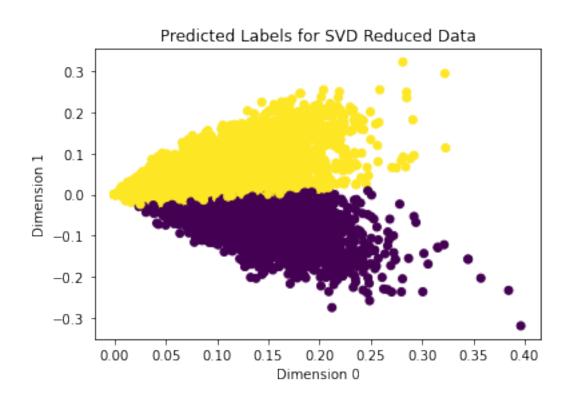
```
[95]: #07
      nmf_best_k = 2
      svd_best_k = 50
      trunc_svd_q3 = TruncatedSVD(n_components=2, random_state=0)
      nmf_model = NMF(n_components=nmf_best_k, init='random', random_state=0,__
       \rightarrowmax_iter=1000)
      nmf_red = nmf_model.fit_transform(X_tfidf)
      nmf_red_SVD = nmf_red#trunc_svd_q3.fit_transform(nmf_red)
      kmeans_clf.fit(nmf_red_SVD)
      nmf_labels = kmeans_clf.labels_
      svd_red = X_svd[:,:svd_best_k]
      svd_red_SVD = trunc_svd_q3.fit_transform(svd_red)
      kmeans_clf.fit(svd_red_SVD)
      svd_labels = kmeans_clf.labels_
      plt.scatter(nmf_red_SVD[:,0], nmf_red_SVD[:,1], c=nmf_labels)
      plt.title("Predicted Labels for NMF Reduced Data")
      plt.xlabel("Dimension 0")
      plt.ylabel("Dimension 1")
```

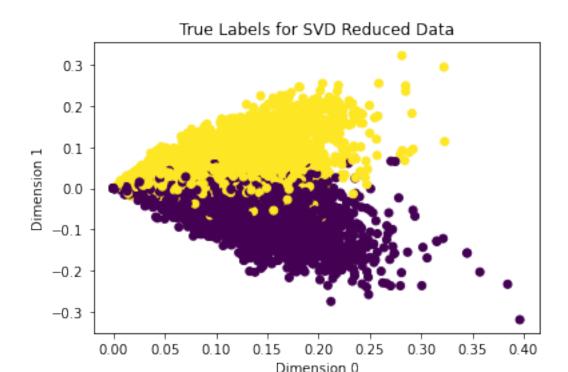
```
plt.show()
plt.scatter(nmf_red_SVD[:,0], nmf_red_SVD[:,1], c=dataset_targets_bin)
plt.title("True Labels for NMF Reduced Data")
plt.xlabel("Dimension 0")
plt.ylabel("Dimension 1")
plt.show()
plt.scatter(svd_red_SVD[:,0], svd_red_SVD[:,1], c=svd_labels)
plt.title("Predicted Labels for SVD Reduced Data")
plt.xlabel("Dimension 0")
plt.ylabel("Dimension 1")
plt.show()
plt.scatter(svd_red_SVD[:,0], svd_red_SVD[:,1], c=dataset_targets_bin)
plt.title("True Labels for SVD Reduced Data")
plt.xlabel("Dimension 0")
plt.ylabel("Dimension 1")
plt.show()
```

Predicted Labels for NMF Reduced Data









```
[96]: #09
      #get all 20 categories this time
     full_dataset = fetch_20newsgroups(subset='all', shuffle = True, random_state = ___
      full_count_vect = CountVectorizer(min_df=3,analyzer=stem_rmv_punc_nolem,_
      ⇔stop_words='english')
     full_X_counts = full_count_vect.fit_transform(full_dataset.data)
     full_X_tfidf = tfidf_transformer.fit_transform(full_X_counts)
     #using SVD with r=50 - best results from earlier
     full_trunc_svd = TruncatedSVD(n_components=1000, random_state=0)
     full_X_svd = full_trunc_svd.fit_transform(full_X_tfidf)
     full_svd_red = full_X_svd[:,:50]
      #k-means w/20 clusters
     full_kmeans_clf = KMeans(n_clusters=20, max_iter=1500, random_state=0, n_init=50)
     full_kmeans_clf.fit(full_svd_red)
      #metrics!
     from scipy.optimize import linear_sum_assignment
     from sklearn.metrics import confusion_matrix
     cm = contingency_matrix(full_dataset.target, full_kmeans_clf.labels_)
```

```
rows, cols=linear_sum_assignment(cm, maximize=True)

plot_mat(cm[rows[:, np.newaxis], cols], xticklabels=cols, yticklabels=rows,__
__size=(15,15), xlabel="Cluster Number", ylabel="Ground Truth Label")

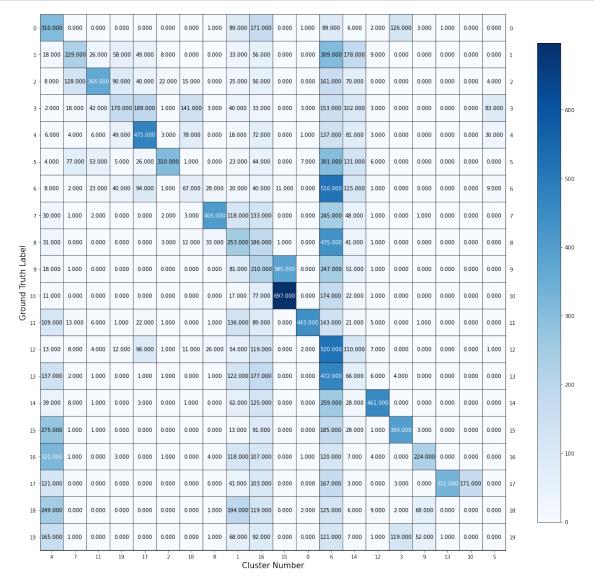
print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target,__
__sfull_kmeans_clf.labels_))

print("Completeness score: ", metrics.completeness_score(full_dataset.target,__
__sfull_kmeans_clf.labels_))

print("V_measure score: ", metrics.v_measure_score(full_dataset.target,__
__sfull_kmeans_clf.labels_))

print("Adjusted Rand Index score: ", metrics.adjusted_rand_score(full_dataset.
__starget, full_kmeans_clf.labels_))

print("Adjusted Mutual Information Score: ",metrics.
__sadjusted_mutual_info_score(full_dataset.target, full_kmeans_clf.labels_))
```



Homogeneity score: 0.31630458145750995 Completeness score: 0.3682552023183464 V_measure score: 0.3403086491478565

Adjusted Rand Index score: 0.10379799750286293

Adjusted Mutual Information Score: 0.3380068278889245

```
[97]: #Q10
      r_{vals} = [2,5,10,30,50,100]
      for i in r_vals:
          nmf_KL = NMF(n_components=i, init='random', random_state=0, max_iter=1000,__
       ⇒beta_loss = "kullback-leibler", solver='mu')
          nmf_red_KL = nmf_KL.fit_transform(full_X_tfidf)
          kmeans_clf_KL = KMeans(n_clusters=20, max_iter=1500, random_state=0,_
       \rightarrown_init=50)
          kmeans_clf_KL.fit(nmf_red_KL)
          print()
          print("r = " + str(i))
          print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target,__
       →kmeans_clf_KL.labels_))
          print("Completeness score: ", metrics.completeness_score(full_dataset.
       →target, kmeans_clf_KL.labels_))
          print("V_measure score: ", metrics.v_measure_score(full_dataset.target,__
       →kmeans_clf_KL.labels_))
          print("Adjusted Rand Index score: ", metrics.
       →adjusted_rand_score(full_dataset.target, kmeans_clf_KL.labels_))
          print("Adjusted Mutual Information Score: ",metrics.
       →adjusted_mutual_info_score(full_dataset.target, kmeans_clf_KL.labels_))
```

r = 2

Homogeneity score: 0.19676527887644293 Completeness score: 0.2136834606423116 V_measure score: 0.2048756966527202

Adjusted Rand Index score: 0.0604569445452653

Adjusted Mutual Information Score: 0.20217480540664073

r = 5

Homogeneity score: 0.35641214262487414 Completeness score: 0.38144347646289517 V_measure score: 0.3685032226345344

Adjusted Rand Index score: 0.15027690007707672

```
r = 10
      Homogeneity score: 0.37714629232162333
      Completeness score: 0.3900505160910544
      V_measure score: 0.3834898799077068
      Adjusted Rand Index score: 0.18833101982676798
      Adjusted Mutual Information Score: 0.38146535004967785
      r = 30
      Homogeneity score: 0.36866650441353743
      Completeness score: 0.42527782891459665
      V_measure score: 0.39495386265506316
      Adjusted Rand Index score: 0.12617792572471453
      Adjusted Mutual Information Score: 0.3928569966801935
      r = 50
      Homogeneity score: 0.30640759897200137
      Completeness score: 0.47873972223954575
      V_measure score: 0.3736610566220568
      Adjusted Rand Index score: 0.05415577061481132
      Adjusted Mutual Information Score: 0.37118024804272814
      r = 100
      Homogeneity score: 0.1924960903730588
      Completeness score: 0.406934904386975
      V_measure score: 0.26135911828245
      Adjusted Rand Index score: 0.028607969901381668
      Adjusted Mutual Information Score: 0.25806962570480285
[153]: #Q11
      import umap
      n_{vals} = [2,5,10,50,100,200,300,500,700,1000]
      metric_scores_euc = []
      metric_scores_cos = []
      for i in n_vals:
          print("working on reducing for n = " + str(i))
          reducer_euc = umap.UMAP(n_components=i, init='random', metric="euclidean")
          embedding_euc = reducer_euc.fit_transform(full_X_tfidf)
          reducer_cos = umap.UMAP(n_components=i, init='random', metric="cosine")
          embedding_cos = reducer_cos.fit_transform(full_X_tfidf)
          print("working on kmeans for n = " + str(i))
```

Adjusted Mutual Information Score: 0.3663744867466823

```
kmeans_clf_euc = KMeans(n_clusters=20, max_iter=1500, random_state=0,__
 \rightarrown_init=50)
    kmeans_clf_euc.fit(embedding_euc)
    kmeans_clf_cos = KMeans(n_clusters=20, max_iter=1500, random_state=0,__
 \rightarrown_init=50)
    kmeans_clf_cos.fit(embedding_cos)
    metric_scores_euc.append((metrics.homogeneity_score(full_dataset.target,_
 →kmeans_clf_euc.labels_),
                               metrics.completeness_score(full_dataset.target,__
 →kmeans_clf_euc.labels_),
                              metrics.v_measure_score(full_dataset.target,__
 →kmeans_clf_euc.labels_),
                              metrics.adjusted_rand_score(full_dataset.target,__
 →kmeans_clf_euc.labels_),
                               metrics.adjusted_mutual_info_score(full_dataset.
 →target, kmeans_clf_euc.labels_)))
    metric_scores_cos.append((metrics.homogeneity_score(full_dataset.target,__
 →kmeans_clf_cos.labels_),
                               metrics.completeness_score(full_dataset.target,__
 →kmeans_clf_cos.labels_),
                              metrics.v_measure_score(full_dataset.target,_
 →kmeans_clf_cos.labels_),
                               metrics.adjusted_rand_score(full_dataset.target,_
 →kmeans_clf_cos.labels_),
                               metrics.adjusted_mutual_info_score(full_dataset.
 →target, kmeans_clf_cos.labels_)))
plt.plot(n_vals, metric_scores_euc)
plt.legend(("Homogeneity", "Completeness", "V_measure", "Adjusted Rand Score", u
 →"Adjusted Mutual Info"))
plt.xlabel("n_components")
plt.ylabel("Score")
plt.title("Scores for Various n_components Using UMAP w/Euclidean Metric")
plt.show()
plt.plot(n_vals, metric_scores_cos)
plt.legend(("Homogeneity", "Completeness", "V_measure", "Adjusted Rand Score", u
→"Adjusted Mutual Info"))
plt.xlabel("n_components")
plt.ylabel("Score")
plt.title("Scores for Various n_components Using UMAP w/Cosine Metric")
plt.show()
```

```
Traceback (most recent call last)
KeyboardInterrupt
<ipython-input-153-abf5f6750690> in <module>()
     11
             print("working on reducing for n = " + str(i))
             reducer_euc = umap.UMAP(n_components=i, init='random',_
 →metric="euclidean")
---> 13
             embedding_euc = reducer_euc.fit_transform(full_X_tfidf)
             reducer_cos = umap.UMAP(n_components=i, init='random', metric="cosine")
     15
             embedding_cos = reducer_cos.fit_transform(full_X_tfidf)
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\umap\umap_.py in__
 →fit_transform(self, X, y)
   2012
                      Embedding of the training data in low-dimensional space.
   2013
-> 2014
                  self.fit(X, y)
   2015
                  return self.embedding_
   2016
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\umap\umap_.py in_
 →fit(self, X, y)
   1799
                           self.low_memory,
   1800
                           use_pynndescent=True,
-> 1801
                           verbose=self.verbose,
   1802
   1803
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\umap\umap_.py in_
 →nearest_neighbors(X, n_neighbors, metric, metric_kwds, angular, random_state, ____
 →low_memory, use_pynndescent, verbose)
    298
                           max_candidates=60,
    299
                           low_memory=low_memory,
--> 300
                           verbose=verbose.
    301
    302
                      knn_indices, knn_dists = nnd.neighbor_graph
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\pynndescent\pynndescent
 →py in __init__(self, data, metric, metric_kwds, n_neighbors, n_trees, u →leaf_size, pruning_degree_multiplier, diversify_prob, n_search_trees, u →tree_init, init_graph, random_state, low_memory, max_candidates, n_iters, u
 →delta, n_jobs, compressed, verbose)
    789
                           current_random_state,
    790
                           self.n_jobs,
--> 791
                           self._angular_trees,
    792
    793
                      leaf_array = rptree_leaf_array(self._rp_forest)
```

```
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\pynndescent\rp_trees.py_
 →in make_forest(data, n_neighbors, n_trees, leaf_size, rng_state, random_state,
 →n_jobs, angular)
    994
                            angular.
    995
                        )
--> 996
                        for i in range(n_trees)
    997
                    )
    998
                else:
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\joblib\parallel.py in_
 →__call__(self, iterable)
   1059
   1060
                    with self._backend.retrieval_context():
-> 1061
                        self.retrieve()
   1062
                    # Make sure that we get a last message telling us we are done
   1063
                    elapsed_time = time.time() - self._start_time
C:\Users\loicm\anaconda3\envs\ece219_2\lib\site-packages\joblib\parallel.py in_
 →retrieve(self)
    938
                    try:
    939
                        if getattr(self._backend, 'supports_timeout', False):
--> 940
                            self._output.extend(job.get(timeout=self.timeout))
    941
                        else:
                            self._output.extend(job.get())
    942
C:\Users\loicm\anaconda3\envs\ece219_2\lib\multiprocessing\pool.py in get(self,_
 →timeout)
    649
    650
            def get(self, timeout=None):
                self.wait(timeout)
--> 651
    652
                if not self.ready():
    653
                    raise TimeoutError
C:\Users\loicm\anaconda3\envs\ece219_2\lib\multiprocessing\pool.py in wait(self,
 →timeout)
    646
    647
            def wait(self, timeout=None):
--> 648
                self._event.wait(timeout)
    649
    650
            def get(self, timeout=None):
C:\Users\loicm\anaconda3\envs\ece219_2\lib\threading.py in wait(self, timeout)
    550
                    signaled = self._flag
    551
                    if not signaled:
--> 552
                        signaled = self._cond.wait(timeout)
    553
                    return signaled
    554
```

```
[143]: #012
       #Best n_components = 5 for both euclidean and cosine metrics
      reducer_euc = umap.UMAP(n_components=5, init='random', metric="euclidean", u
       →random_state=42)
      embedding_euc = reducer_euc.fit_transform(full_X_tfidf)
      reducer_cos = umap.UMAP(n_components=5, init='random', metric="cosine", __
       →random_state=42)
      embedding_cos = reducer_cos.fit_transform(full_X_tfidf)
      kmeans_clf_euc = KMeans(n_clusters=20, max_iter=1500, random_state=0, n_init=50)
      kmeans_clf_euc.fit(embedding_euc)
      kmeans_clf_cos = KMeans(n_clusters=20, max_iter=1500, random_state=0, n_init=50)
      kmeans_clf_cos.fit(embedding_cos)
      print("Performance of UMAP reduced data for n_components=5")
      print()
      print("Euclidean metric")
      print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target, u
       →kmeans_clf_euc.labels_))
      print("Completeness score: ", metrics.completeness_score(full_dataset.target,__
        →kmeans_clf_euc.labels_))
      print("V_measure score: ", metrics.v_measure_score(full_dataset.target,__
        →kmeans_clf_euc.labels_))
      print("Adjusted Rand Index score: ", metrics.adjusted_rand_score(full_dataset.
       →target, kmeans_clf_euc.labels_))
      print("Adjusted Mutual Information Score: ",metrics.
        -adjusted_mutual_info_score(full_dataset.target, kmeans_clf_euc.labels_))
      print()
      print("Cosine metric")
      print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target,_
       →kmeans_clf_cos.labels_))
      print("Completeness score: ", metrics.completeness_score(full_dataset.target,_
        →kmeans_clf_cos.labels_))
      print("V_measure score: ", metrics.v_measure_score(full_dataset.target,_
        →kmeans_clf_cos.labels_))
```

Performance of UMAP reduced data for n_components=5

Euclidean metric

Homogeneity score: 0.008428056566878686 Completeness score: 0.008490790660743977 V_measure score: 0.008459307306639803

Adjusted Rand Index score: 0.0014762057832842827

Adjusted Mutual Information Score: 0.00524781170211466

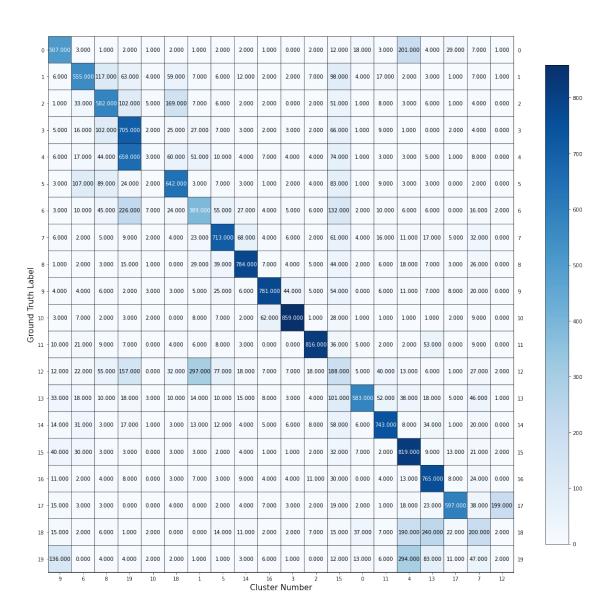
Cosine metric

Homogeneity score: 0.5347926404764849 Completeness score: 0.5549487009354954 V_measure score: 0.5446842655666289

Adjusted Rand Index score: 0.4167210932677909

Adjusted Mutual Information Score: 0.5431756896373161

0	49.000	35.000	30.000	28.000	37.000	38.000	29.000	35.000	29.000	44.000	83.000	45.000	64.000	33.000	35.000	44.000	31.000	33.000	37.000	40.000	0	
1	- 39.000	56.000	52.000	39.000	46.000	54.000	60.000	39.000	40.000	44.000	88.000	41.000	70.000	45.000	43.000	59.000	49.000	24.000	40.000	45.000	1	
	- 34.000	30.000		33.000	33.000		44.000	36.000	47.000	49.000	103.000		80.000	38.000	47.000	41.000		59.000	45.000	24.000	2	- 100
	- 39.000	44.000	38.000	74.000	57.000	46.000	55.000	35.000	36.000	61.000	91.000	65.000	72.000	46.000	51.000	51.000	35.000	22.000	17.000	47.000	3	
4	22.000	42.000	56.000	49.000	63.000	50.000	54.000	47.000	39.000	56.000	66.000	22.000	79.000	41.000	53.000	56.000	52.000	35.000	37.000	44.000	4	
5 -	42.000	47.000	35.000	41.000	54.000	56.000	58.000	34.000	40.000	49.000	80.000	35.000	77.000	50.000	61.000	55.000	56.000	32.000	41.000	45.000	5	
6	- 39.000	47.000	46.000	51.000	59.000	50.000	72.000	40.000	34.000	64.000	84.000	25.000	75.000	64.000	48.000	53.000	49.000	12.000	13.000	50.000	6	- 80
7	- 34.000	35.000	39.000	32.000	45.000	61.000	43.000	47.000	38.000	60.000	103.000	47.000	68.000	39.000	58.000	50.000	61.000	50.000	33.000	47.000	7	
	32.000	41.000	45.000	36.000	49.000	50.000	44.000	34.000	44.000	62.000	81.000	58.000	77.000	46.000	58.000	56.000	55.000	43.000	44.000	41.000	8	
Ground Truth Label	- 35.000	42.000	55.000	38.000	37.000	43.000	70.000	32.000	48.000	73.000	95.000	58.000	80.000	32.000	52.000	53.000	36.000	58.000	14.000	43.000	9	
od Trut	19.000	41.000	35.000	40.000	52.000	51.000	56.000	40.000	42.000	47.000	108.000	81.000	75.000	47.000	44.000	50.000	33.000	70.000	42.000	26.000	10	- 60
Broul 31	43.000	39.000	44.000	32.000	50.000	60.000	40.000	23.000	46.000	61.000	85.000	82.000	83.000	35.000	36.000	63.000	39.000	29.000	64.000	37.000	11	
12	- 38.000	35 000	51 000	37 000	46 000	47.000	42 000	40 000	44 000	49 000	97.000	44 000	82 000	37 000	54 000	52 000	44 000	39.000	65 000	41.000	12	
	45.000							39.000							55.000			50.000		45.000		
13	45.000	54.000	44.000	26.000	42.000	47.000	46.000	39.000	35.000	63.000	71.000	57.000	68.000	52.000	35.000	45.000	35.000	50.000	49.000	45.000	13	- 40
14	- 31.000	41.000	50.000	41.000	47.000	54.000	45.000	38.000	36.000	62.000	91.000	62.000	77.000	36.000	61.000	51.000	45.000	34.000	37.000	48.000	14	
15	43.000	52.000	52.000	37.000	49.000	47.000	40.000	54.000	54.000			38.000		33.000	54.000		53.000	24.000	26.000		15	
16	31.000	35.000	36.000	34.000	58.000	35.000	34.000	35.000	27.000	55.000		62.000	81.000	34.000	41.000	42.000	62.000	55.000	45.000	34.000	16	
17	25.000	34.000	46.000	22.000	34.000	30.000	40.000	16.000	35.000	44.000	81.000	81.000	65.000	43.000	51.000	47.000	51.000	70.000	92.000	33.000	17	- 20
18	15.000	28.000	39.000	21.000	47.000	37.000	37.000	32.000	29.000	37.000	58.000	28.000	61.000	34.000	38.000	33.000	37.000	50.000	89.000	25.000	18	
19	11.000	24.000	27.000	25.000	23.000	34.000	24.000	24.000	19.000	41.000	55.000	33.000	41.000	25.000	41.000	29.000	25.000	49.000	37.000	41.000	19	
	18	8	14	12	7	6	i	10	16 CI	uster	15 Numb	er	19	2	17	9	4	3	5	13		



```
[142]: #Q13
#using same umap parameters as above

from sklearn.cluster import AgglomerativeClustering

umap_best_n = 5
umap_best_metric = 'cosine'

umap_reducer = umap.UMAP(n_components=umap_best_n, init='random', umap_reducer.get_n)

ometric=umap_best_metric, random_state=42)

red_umap = umap_reducer.fit_transform(full_X_tfidf)
```

```
agg_clust_ward = AgglomerativeClustering(n_clusters=20, linkage='ward')
agg_clust_single = AgglomerativeClustering(n_clusters=20, linkage='single')
agg_clust_ward.fit(red_umap)
agg_clust_single.fit(red_umap)
print("Ward Linkage Scores")
print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target,__
 →agg_clust_ward.labels_))
print("Completeness score: ", metrics.completeness_score(full_dataset.target, __
 →agg_clust_ward.labels_))
print("V_measure score: ", metrics.v_measure_score(full_dataset.target,_
 →agg_clust_ward.labels_))
print("Adjusted Rand Index score: ", metrics.adjusted_rand_score(full_dataset.
 →target, agg_clust_ward.labels_))
print("Adjusted Mutual Information Score: ",metrics.
 →adjusted_mutual_info_score(full_dataset.target, agg_clust_ward.labels_))
print()
print("Single Linkage Scores")
print("Homogeneity score: ", metrics.homogeneity_score(full_dataset.target,_
 →agg_clust_single.labels_))
print("Completeness score: ", metrics.completeness_score(full_dataset.target,_
 →agg_clust_single.labels_))
print("V_measure score: ", metrics.v_measure_score(full_dataset.target,_
 →agg_clust_single.labels_))
print("Adjusted Rand Index score: ", metrics.adjusted_rand_score(full_dataset.
 →target, agg_clust_single.labels_))
print("Adjusted Mutual Information Score: ",metrics.
 -adjusted_mutual_info_score(full_dataset.target, agg_clust_single.labels_))
Ward Linkage Scores
Homogeneity score: 0.524473300498607
Completeness score: 0.5479912976714623
V_measure score: 0.53597443687123
Adjusted Rand Index score: 0.3940501791969448
Adjusted Mutual Information Score: 0.5344316766896055
Single Linkage Scores
Homogeneity score: 0.022917457355135236
Completeness score: 0.3464156710750703
V_measure score: 0.04299081646294147
Adjusted Rand Index score: 0.00046761208355549543
Adjusted Mutual Information Score: 0.03748616963750802
```

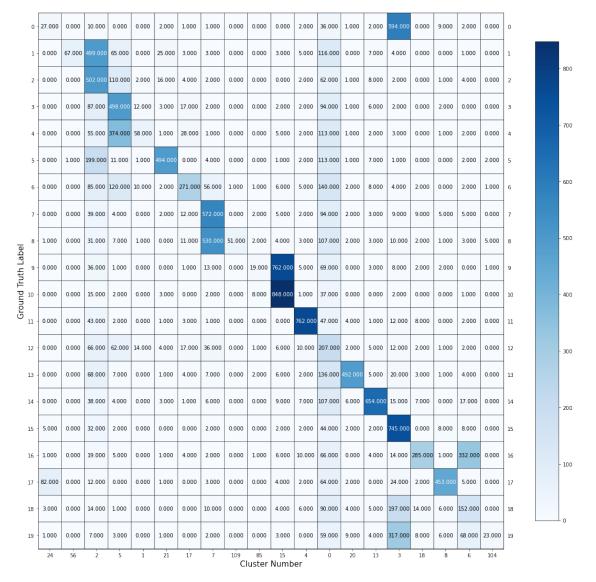
```
[140]: #Q14 with more hyperparams
      eps_vals = [.01, .05, .1, .3, .5, .7, .9, 1]
      metric_types = [ 'euclidean', 'manhattan', 'cosine']
      min_samples = [10, 20, 50, 70, 100]
      db_max_score = -100
      hdb_max_score = -100
      for eps in eps_vals:
          for metric_t in metric_types:
               for min_samp in min_samples:
                   print("Current Iteration: eps={:f}, metric={:s}, min_samples={:d}".
        →format(eps, metric_t, min_samp))
                   DBSCAN_clust = DBSCAN(min_samples=min_samp, metric=metric_t, eps=eps)
                   #HDB does not support cosine metric
                   if(metric_t != 'cosine'):
                       HDBSCAN_clust = HDBSCAN(min_cluster_size=100,__
        →min_samples=min_samp, cluster_selection_epsilon=eps)
                   DBSCAN_clust.fit(red_umap)
                   HDBSCAN_clust.fit(red_umap)
                   homog_score = metrics.homogeneity_score(full_dataset.target,_
        →DBSCAN_clust.labels_)
                   comp_score = metrics.completeness_score(full_dataset.target,___
        →DBSCAN_clust.labels_)
                   v_score = metrics.v_measure_score(full_dataset.target, DBSCAN_clust.
        →labels_)
                   rand_score = metrics.adjusted_rand_score(full_dataset.target,_
        →DBSCAN_clust.labels_)
                   adj_mut_score = metrics.adjusted_mutual_info_score(full_dataset.
        →target, DBSCAN_clust.labels_)
                   score_avg = (homog_score + comp_score + v_score + rand_score +
        →adj_mut_score)/5
                   labels = DBSCAN_clust.labels_
                   n_clust = len(set(labels)) - (1 if -1 in labels else 0)
                   if score_avg > db_max_score:
                       db_max_score = score_avg
                       db_max_params = (eps, metric_t, min_samp)
                       db_max_metrics = (homog_score, comp_score, v_score, rand_score,
        →adj_mut_score)
```

```
homog_score = metrics.homogeneity_score(full_dataset.target,_
 →HDBSCAN_clust.labels_)
            comp_score = metrics.completeness_score(full_dataset.target,__
 →HDBSCAN_clust.labels_)
            v_score = metrics.v_measure_score(full_dataset.target, HDBSCAN_clust.
 →labels_)
            rand_score = metrics.adjusted_rand_score(full_dataset.target,_
 →HDBSCAN_clust.labels_)
            adj_mut_score = metrics.adjusted_mutual_info_score(full_dataset.
 →target, HDBSCAN_clust.labels_)
            score_avg = (homog_score + comp_score + v_score + rand_score +
 →adj_mut_score)/5
            labels = HDBSCAN_clust.labels_
            n_clust = len(set(labels)) - (1 if -1 in labels else 0)
            if metric_t != 'cosine' and score_avg > hdb_max_score:
                hdb_max_score = score_avg
                hdb_max_params = (eps, metric_t, min_samp)
                hdb_max_metrics = (homog_score, comp_score, v_score, rand_score,_
 →adj_mut_score)
print()
print("DBSCAN Best Parameters: eps={:f}, metric={:s}, min_samples={:d}".
 format(db_max_params[0], db_max_params[1], db_max_params[2]))
print()
print("Best DBSCAN Scores")
print("Homogeneity score: ", db_max_metrics[0])
print("Completeness score: ", db_max_metrics[1])
print("V_measure score: ", db_max_metrics[2])
print("Adjusted Rand Index score: ", db_max_metrics[3])
print("Adjusted Mutual Information Score: ", db_max_metrics[4])
print()
print("HDBSCAN Best Parameters: eps={:f}, metric={:s}, min_samples={:d}".
 →format(hdb_max_params[0], hdb_max_params[1], hdb_max_params[2]))
print()
print("Best HDBSCAN Scores")
print("Homogeneity score: ", hdb_max_metrics[0])
print("Completeness score: ", hdb_max_metrics[1])
print("V_measure score: ", hdb_max_metrics[2])
print("Adjusted Rand Index score: ", hdb_max_metrics[3])
print("Adjusted Mutual Information Score: ", hdb_max_metrics[4])
```

```
Current Iteration: eps=0.010000, metric=euclidean, min_samples=10
Current Iteration: eps=0.010000, metric=euclidean, min_samples=20
Current Iteration: eps=0.010000, metric=euclidean, min_samples=50
Current Iteration: eps=0.010000, metric=euclidean, min_samples=70
Current Iteration: eps=0.010000, metric=euclidean, min_samples=100
Current Iteration: eps=0.010000, metric=manhattan, min_samples=10
Current Iteration: eps=0.010000, metric=manhattan, min_samples=20
Current Iteration: eps=0.010000, metric=manhattan, min_samples=50
Current Iteration: eps=0.010000, metric=manhattan, min_samples=70
Current Iteration: eps=0.010000, metric=manhattan, min_samples=100
Current Iteration: eps=0.010000, metric=cosine, min_samples=10
Current Iteration: eps=0.010000, metric=cosine, min_samples=20
Current Iteration: eps=0.010000, metric=cosine, min_samples=50
Current Iteration: eps=0.010000, metric=cosine, min_samples=70
Current Iteration: eps=0.010000, metric=cosine, min_samples=100
Current Iteration: eps=0.050000, metric=euclidean, min_samples=10
Current Iteration: eps=0.050000, metric=euclidean, min_samples=20
Current Iteration: eps=0.050000, metric=euclidean, min_samples=50
Current Iteration: eps=0.050000, metric=euclidean, min_samples=70
Current Iteration: eps=0.050000, metric=euclidean, min_samples=100
Current Iteration: eps=0.050000, metric=manhattan, min_samples=10
Current Iteration: eps=0.050000, metric=manhattan, min_samples=20
Current Iteration: eps=0.050000, metric=manhattan, min_samples=50
Current Iteration: eps=0.050000, metric=manhattan, min_samples=70
Current Iteration: eps=0.050000, metric=manhattan, min_samples=100
Current Iteration: eps=0.050000, metric=cosine, min_samples=10
Current Iteration: eps=0.050000, metric=cosine, min_samples=20
Current Iteration: eps=0.050000, metric=cosine, min_samples=50
Current Iteration: eps=0.050000, metric=cosine, min_samples=70
Current Iteration: eps=0.050000, metric=cosine, min_samples=100
Current Iteration: eps=0.100000, metric=euclidean, min_samples=10
Current Iteration: eps=0.100000, metric=euclidean, min_samples=20
Current Iteration: eps=0.100000, metric=euclidean, min_samples=50
Current Iteration: eps=0.100000, metric=euclidean, min_samples=70
Current Iteration: eps=0.100000, metric=euclidean, min_samples=100
Current Iteration: eps=0.100000, metric=manhattan, min_samples=10
Current Iteration: eps=0.100000, metric=manhattan, min_samples=20
Current Iteration: eps=0.100000, metric=manhattan, min_samples=50
Current Iteration: eps=0.100000, metric=manhattan, min_samples=70
Current Iteration: eps=0.100000, metric=manhattan, min_samples=100
Current Iteration: eps=0.100000, metric=cosine, min_samples=10
Current Iteration: eps=0.100000, metric=cosine, min_samples=20
Current Iteration: eps=0.100000, metric=cosine, min_samples=50
Current Iteration: eps=0.100000, metric=cosine, min_samples=70
Current Iteration: eps=0.100000, metric=cosine, min_samples=100
Current Iteration: eps=0.300000, metric=euclidean, min_samples=10
Current Iteration: eps=0.300000, metric=euclidean, min_samples=20
Current Iteration: eps=0.300000, metric=euclidean, min_samples=50
```

```
Current Iteration: eps=0.300000, metric=euclidean, min_samples=70
Current Iteration: eps=0.300000, metric=euclidean, min_samples=100
Current Iteration: eps=0.300000, metric=manhattan, min_samples=10
Current Iteration: eps=0.300000, metric=manhattan, min_samples=20
Current Iteration: eps=0.300000, metric=manhattan, min_samples=50
Current Iteration: eps=0.300000, metric=manhattan, min_samples=70
Current Iteration: eps=0.300000, metric=manhattan, min_samples=100
Current Iteration: eps=0.300000, metric=cosine, min_samples=10
Current Iteration: eps=0.300000, metric=cosine, min_samples=20
Current Iteration: eps=0.300000, metric=cosine, min_samples=50
Current Iteration: eps=0.300000, metric=cosine, min_samples=70
Current Iteration: eps=0.300000, metric=cosine, min_samples=100
Current Iteration: eps=0.500000, metric=euclidean, min_samples=10
Current Iteration: eps=0.500000, metric=euclidean, min_samples=20
Current Iteration: eps=0.500000, metric=euclidean, min_samples=50
Current Iteration: eps=0.500000, metric=euclidean, min_samples=70
Current Iteration: eps=0.500000, metric=euclidean, min_samples=100
Current Iteration: eps=0.500000, metric=manhattan, min_samples=10
Current Iteration: eps=0.500000, metric=manhattan, min_samples=20
Current Iteration: eps=0.500000, metric=manhattan, min_samples=50
Current Iteration: eps=0.500000, metric=manhattan, min_samples=70
Current Iteration: eps=0.500000, metric=manhattan, min_samples=100
Current Iteration: eps=0.500000, metric=cosine, min_samples=10
Current Iteration: eps=0.500000, metric=cosine, min_samples=20
Current Iteration: eps=0.500000, metric=cosine, min_samples=50
Current Iteration: eps=0.500000, metric=cosine, min_samples=70
Current Iteration: eps=0.500000, metric=cosine, min_samples=100
Current Iteration: eps=0.700000, metric=euclidean, min_samples=10
Current Iteration: eps=0.700000, metric=euclidean, min_samples=20
Current Iteration: eps=0.700000, metric=euclidean, min_samples=50
Current Iteration: eps=0.700000, metric=euclidean, min_samples=70
Current Iteration: eps=0.700000, metric=euclidean, min_samples=100
Current Iteration: eps=0.700000, metric=manhattan, min_samples=10
Current Iteration: eps=0.700000, metric=manhattan, min_samples=20
Current Iteration: eps=0.700000, metric=manhattan, min_samples=50
Current Iteration: eps=0.700000, metric=manhattan, min_samples=70
Current Iteration: eps=0.700000, metric=manhattan, min_samples=100
Current Iteration: eps=0.700000, metric=cosine, min_samples=10
Current Iteration: eps=0.700000, metric=cosine, min_samples=20
Current Iteration: eps=0.700000, metric=cosine, min_samples=50
Current Iteration: eps=0.700000, metric=cosine, min_samples=70
Current Iteration: eps=0.700000, metric=cosine, min_samples=100
Current Iteration: eps=0.900000, metric=euclidean, min_samples=10
Current Iteration: eps=0.900000, metric=euclidean, min_samples=20
Current Iteration: eps=0.900000, metric=euclidean, min_samples=50
Current Iteration: eps=0.900000, metric=euclidean, min_samples=70
Current Iteration: eps=0.900000, metric=euclidean, min_samples=100
Current Iteration: eps=0.900000, metric=manhattan, min_samples=10
```

```
Current Iteration: eps=0.900000, metric=manhattan, min_samples=20
      Current Iteration: eps=0.900000, metric=manhattan, min_samples=50
      Current Iteration: eps=0.900000, metric=manhattan, min_samples=70
      Current Iteration: eps=0.900000, metric=manhattan, min_samples=100
      Current Iteration: eps=0.900000, metric=cosine, min_samples=10
      Current Iteration: eps=0.900000, metric=cosine, min_samples=20
      Current Iteration: eps=0.900000, metric=cosine, min_samples=50
      Current Iteration: eps=0.900000, metric=cosine, min_samples=70
      Current Iteration: eps=0.900000, metric=cosine, min_samples=100
      Current Iteration: eps=1.000000, metric=euclidean, min_samples=10
      Current Iteration: eps=1.000000, metric=euclidean, min_samples=20
      Current Iteration: eps=1.000000, metric=euclidean, min_samples=50
      Current Iteration: eps=1.000000, metric=euclidean, min_samples=70
      Current Iteration: eps=1.000000, metric=euclidean, min_samples=100
      Current Iteration: eps=1.000000, metric=manhattan, min_samples=10
      Current Iteration: eps=1.000000, metric=manhattan, min_samples=20
      Current Iteration: eps=1.000000, metric=manhattan, min_samples=50
      Current Iteration: eps=1.000000, metric=manhattan, min_samples=70
      Current Iteration: eps=1.000000, metric=manhattan, min_samples=100
      Current Iteration: eps=1.000000, metric=cosine, min_samples=10
      Current Iteration: eps=1.000000, metric=cosine, min_samples=20
      Current Iteration: eps=1.000000, metric=cosine, min_samples=50
      Current Iteration: eps=1.000000, metric=cosine, min_samples=70
      Current Iteration: eps=1.000000, metric=cosine, min_samples=100
      DBSCAN Best Parameters: eps=0.500000, metric=manhattan, min_samples=10
      Best DBSCAN Scores
      Homogeneity score: 0.5407607595568981
      Completeness score: 0.45471096739764455
      V_measure score: 0.49401673890037173
      Adjusted Rand Index score: 0.2967260971944622
      Adjusted Mutual Information Score: 0.47920988791262803
      HDBSCAN Best Parameters: eps=0.300000, metric=euclidean, min_samples=10
      Best HDBSCAN Scores
      Homogeneity score: 0.43889528158878366
      Completeness score: 0.5288819318072416
      V_measure score: 0.47970500064413385
      Adjusted Rand Index score: 0.2565568616115164
      Adjusted Mutual Information Score: 0.4780526456871693
[141]: # Q15
      DBSCAN_clust = DBSCAN(min_samples=10, eps=0.5, metric='manhattan')
      DBSCAN_clust.fit(red_umap)
```



Estimated number of clusters: 168

```
[126]: #016
       #get data and vectorize
       #Using the larger training dataset
      import pandas
      from sklearn.cluster import DBSCAN
      from sklearn.metrics.cluster import contingency_matrix
      from scipy.optimize import linear_sum_assignment
      from plotmat import plot_mat
      import numpy as np
      import umap.umap_ as umap
      bbc = pandas.read_csv('BBC_News_Train.csv')
      categories = (bbc.Category.unique()).tolist()
      targets = []
      for i in bbc.Category:
          for j in range(0,len(categories)):
               if i == categories[j]:
                   targets.append(j)
      bbc['target'] = targets
      X_counts = count_vect.fit_transform(bbc.Text)
      bbc_X_tfidf = tfidf_transformer.fit_transform(X_counts)
      print("bbc_X_tfidf shape: ", bbc_X_tfidf.shape)
```

bbc_X_tfidf shape: (1490, 10197)

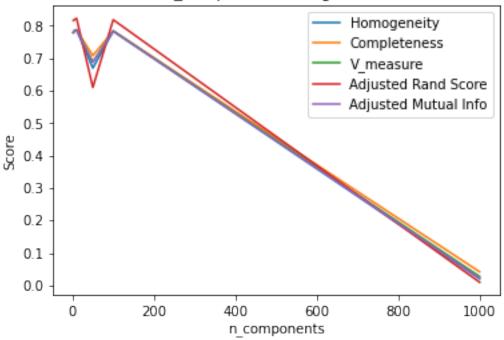
```
col_mean = np.nanmean(embedding_cos, axis=0)
    inds = np.where(np.isnan(embedding_cos))
    embedding_cos[inds] = np.take(col_mean, inds[1])
    print("working on kmeans for n = " + str(i))
    kmeans_clf_euc = KMeans(n_clusters=5, max_iter=1500, random_state=0,_
 \rightarrown_init=50)
    kmeans_clf_euc.fit(embedding_euc)
    kmeans_clf_cos = KMeans(n_clusters=5, max_iter=1500, random_state=0,_u
 \rightarrown_init=50)
    kmeans_clf_cos.fit(embedding_cos)
    metric_scores_euc.append((metrics.homogeneity_score(bbc.target,_
 →kmeans_clf_euc.labels_),
                               metrics.completeness_score(bbc.target,_
 →kmeans_clf_euc.labels_),
                              metrics.v_measure_score(bbc.target, kmeans_clf_euc.
 →labels_),
                               metrics.adjusted_rand_score(bbc.target,__
 →kmeans_clf_euc.labels_),
                               metrics.adjusted_mutual_info_score(bbc.target,__
 →kmeans_clf_euc.labels_)))
    metric_scores_cos.append((metrics.homogeneity_score(bbc.target,_
 →kmeans_clf_cos.labels_),
                               metrics.completeness_score(bbc.target,_
 →kmeans_clf_cos.labels_),
                              metrics.v_measure_score(bbc.target, kmeans_clf_cos.
 →labels_),
                              metrics.adjusted_rand_score(bbc.target,__
 →kmeans_clf_cos.labels_),
                              metrics.adjusted_mutual_info_score(bbc.target,__
 →kmeans_clf_cos.labels_)))
# Plot Metrics against n_components
plt.plot(n_vals, metric_scores_euc)
plt.legend(("Homogeneity", "Completeness", "V_measure", "Adjusted Rand Score", __
→"Adjusted Mutual Info"))
plt.xlabel("n_components")
plt.ylabel("Score")
plt.title("Scores for Various n_components Using UMAP w/Euclidean Metric")
plt.show()
plt.plot(n_vals, metric_scores_cos)
```

```
plt.legend(("Homogeneity", "Completeness", "V_measure", "Adjusted Rand Score", u
 →"Adjusted Mutual Info"))
plt.xlabel("n_components")
plt.ylabel("Score")
plt.title("Scores for Various n_components Using UMAP w/Cosine Metric")
plt.show()
# Find best values and print metrics
avs_Kmeans_cos = []
for i in range(len(metric_scores_cos)):
 →((metric_scores_cos[i][0]+metric_scores_cos[i][1]+metric_scores_cos[i][2]+metrid_scores_cos[i]
 →5)
    avs_Kmeans_cos.append(a)
avs_Kmeans_euc = []
for i in range(len(metric_scores_euc)):
 →((metric_scores_euc[i][0]+metric_scores_euc[i][1]+metric_scores_euc[i][2]+metric_scores_euc[i
    avs_Kmeans_euc.append(a)
print('Best Kmeans Cos: ', max(avs_Kmeans_cos))
print('Best Kmeans Euc: ', max(avs_Kmeans_euc))
c = avs_Kmeans_euc.index(max(avs_Kmeans_euc))
b = avs_Kmeans_cos.index(max(avs_Kmeans_cos))
print()
print("Best KMeans EUC Scores")
print("Homogeneity score: ", metric_scores_euc[c][0])
print("Completeness score: ", metric_scores_euc[c][1])
print("V_measure score: ", metric_scores_euc[c][2])
print("Adjusted Rand Index score: ", metric_scores_euc[c][3])
print("Adjusted Mutual Information Score: ", metric_scores_euc[c][4])
print("Best N_Components: ", n_vals[c])
umap_reducer = umap.UMAP(n_components = n_vals[c], init = 'random', metric = __

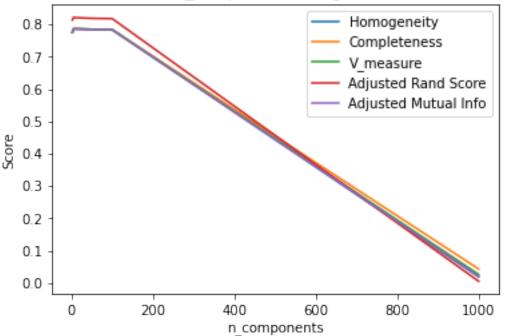
→ 'euclidean', random_state=0)
bbc_umap = umap_reducer.fit_transform(bbc_X_tfidf)
#k-means w/5 clusters
bbc_kmeans_clf = KMeans(n_clusters=5, max_iter=1500, random_state=0, n_init=50)
bbc_kmeans_clf.fit(bbc_umap)
#metrics!
```

```
from scipy.optimize import linear_sum_assignment
from sklearn.metrics import confusion_matrix
cm = contingency_matrix(bbc.target, bbc_kmeans_clf.labels_)
rows, cols=linear_sum_assignment(cm, maximize=True)
plot_mat(cm[rows[:, np.newaxis], cols], xticklabels=cols, yticklabels=rows,_u
 \rightarrowsize=(8,8))
print()
print("Best KMeans COS Scores")
print("Homogeneity score: ", metric_scores_cos[b][0])
print("Completeness score: ", metric_scores_cos[b][1])
print("V_measure score: ", metric_scores_cos[b][2])
print("Adjusted Rand Index score: ", metric_scores_cos[b][3])
print("Adjusted Mutual Information Score: ", metric_scores_cos[b][4])
print("Best N_Components: ", n_vals[b])
umap_reducer = umap.UMAP(n_components = n_vals[c], init = 'random', metric = __
 bbc_umap = umap_reducer.fit_transform(bbc_X_tfidf)
#k-means w/5 clusters
bbc_kmeans_clf = KMeans(n_clusters=5, max_iter=1500, random_state=0, n_init=50)
bbc_kmeans_clf.fit(bbc_umap)
#metrics!
from scipy.optimize import linear_sum_assignment
from sklearn.metrics import confusion_matrix
cm = contingency_matrix(bbc.target, bbc_kmeans_clf.labels_)
rows, cols=linear_sum_assignment(cm, maximize=True)
plot_mat(cm[rows[:, np.newaxis], cols], xticklabels=cols, yticklabels=rows,_u
 \rightarrowsize=(8,8))
working on reducing for n = 2
working on kmeans for n = 2
working on reducing for n = 5
working on kmeans for n = 5
working on reducing for n = 10
working on kmeans for n = 10
working on reducing for n = 50
working on kmeans for n = 50
working on reducing for n = 100
working on kmeans for n = 100
working on reducing for n = 1000
working on kmeans for n = 1000
```

Scores for Various n_components Using UMAP w/Euclidean Metric



Scores for Various n_components Using UMAP w/Cosine Metric



Best Kmeans Cos: 0.7930692963971817 Best Kmeans Euc: 0.7944233592691203

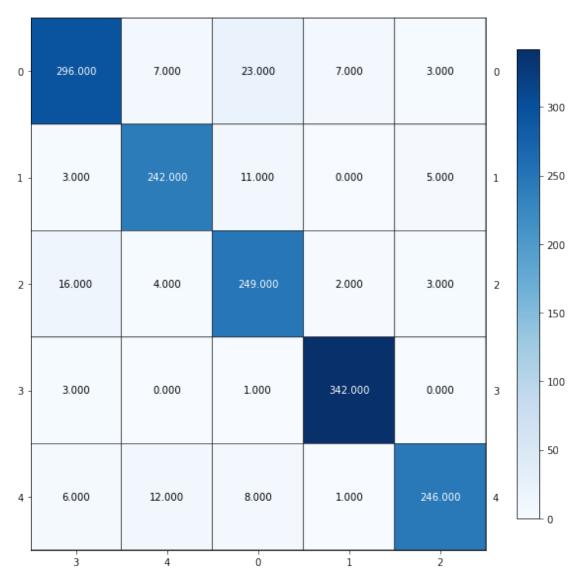
Best KMeans EUC Scores

Homogeneity score: 0.7874360944663693 Completeness score: 0.7874929063651126 V_measure score: 0.7874644993910623

Adjusted Rand Index score: 0.8229763601775667

Adjusted Mutual Information Score: 0.786746935945491

Best N_Components: 10



Best KMeans COS Scores

Homogeneity score: 0.7863911146778318

Completeness score: 0.7863053396249526 V_measure score: 0.7863482248123015

Adjusted Rand Index score: 0.8206748429226303

Adjusted Mutual Information Score: 0.785626959948192

Best N_Components: 10



```
[151]: # Find best parameter combo in DBSCAN with COSINE metric
# Generate 3D surface plots for each metric against epsillon and min samples
\[ \text{-values} \]

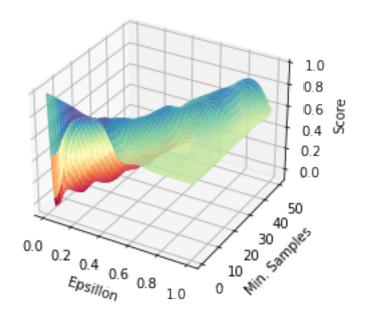
umap_reducer = umap.UMAP(n_components = 5, init = 'random', metric = 'cosine', \[ \text{-random_state=0} \]

bbc_umap_COS = umap_reducer.fit_transform(bbc_X_tfidf)
```

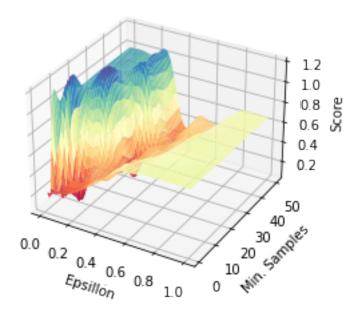
```
eps_vals = [.01, .05, .1, .2, .3, .4, .5, .6, .7, .8, .9, 1]
samples = [1,3,5,10,20,30,35,40,50]
metric_scores_DBSCAN_COS = []
for eps in eps_vals:
    for sam in samples:
        DBSCAN_clust = DBSCAN(min_samples=sam, eps=eps)
        DBSCAN_clust.fit(bbc_umap)
        labels = DBSCAN_clust.labels_
        n_clusters_ = len(set(labels)) - (1 if -1 in labels else 0)
        metric_scores_DBSCAN_COS.append([metrics.homogeneity_score(bbc.target,_
 →DBSCAN_clust.labels_),
                                     metrics.completeness_score(bbc.target,__
 →DBSCAN_clust.labels_),
                                     metrics.v_measure_score(bbc.target,__
 →DBSCAN_clust.labels_),
                                     metrics.adjusted_rand_score(bbc.target,_
 →DBSCAN_clust.labels_),
                                     metrics.adjusted_mutual_info_score(bbc.
→target, DBSCAN_clust.labels_),
                                     n_clusters_,eps,sam])
# Visualize parameter effects on metrics
from mpl_toolkits.mplot3d import Axes3D
from scipy.interpolate import griddata
metricdata = np.array(metric_scores_DBSCAN_COS)
metric_scores = ['Homogeneity', 'Completeness', 'V Measure', 'Adjusted Rand', |
→'Adj. Mutual Info', 'Number of Clusters']
for i in range(0,6):
   plotdata = metricdata[:, [i,6,7]]
    z, x, y = zip(*plotdata)
    z = list(map(float, z))
    grid_x, grid_y = np.mgrid[min(x):max(x):100j, min(y):max(y):100j]
    grid_z = griddata((x, y), z, (grid_x, grid_y), method='cubic')
   fig = plt.figure()
    ax = fig.gca(projection='3d')
    ax.plot_surface(grid_x, grid_y, grid_z, cmap=plt.cm.Spectral)
    ax.set_xlabel('Epsillon')
    ax.set_ylabel('Min. Samples')
```

```
ax.set_zlabel('Score')
    ax.set_title(str(metric_scores[i]))
    plt.show()
# Find best parameter combination for epsillon and minimum samples
avs = []
for i in range(len(metricdata)):
→((metricdata[i][0]+metricdata[i][1]+metricdata[i][2]+metricdata[i][3]+metricdata[i][4])/
 →5)
    avs.append(a)
b = avs.index(max(avs))
print('Best Epsillon = ', metricdata[b][6])
print('Best N_Samples = ', metricdata[b][7])
print()
cm = contingency_matrix(bbc.target, DBSCAN_clust.labels_)
rows, cols=linear_sum_assignment(cm, maximize=True)
plot_mat(cm[rows[:, np.newaxis], cols], xticklabels=cols, yticklabels=rows,_u
\rightarrowsize=(5,5))
print("Estimated Number of Clusters: ", metricdata[b][5])
print()
print("Best DBSCAN Scores for Cosine Distance Metric")
print("Homogeneity score: ", metricdata[b][0])
print("Completeness score: ", metricdata[b][1])
print("V_measure score: ", metricdata[b][2])
print("Adjusted Rand Index score: ", metricdata[b][3])
print("Adjusted Mutual Information Score: ", metricdata[b][4])
```

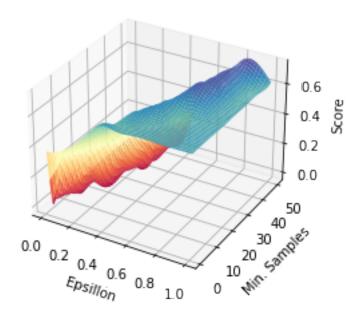
Homogeneity



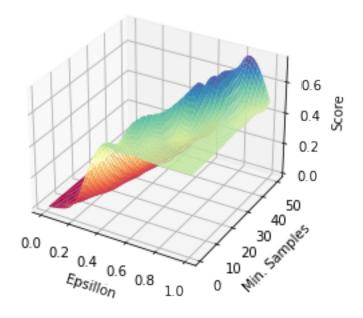
Completeness



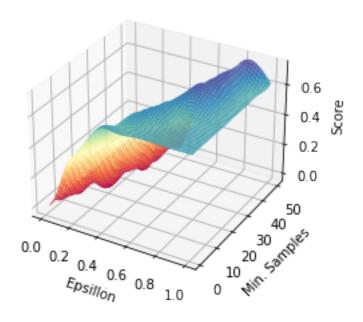
V Measure



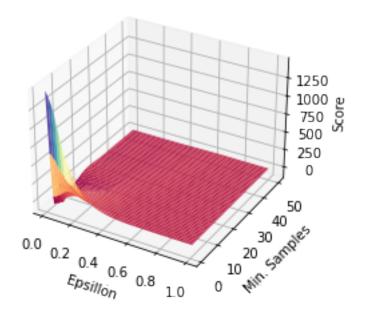
Adjusted Rand



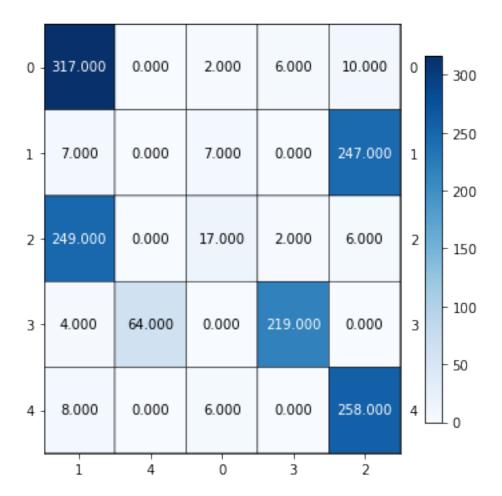
Adj. Mutual Info



Number of Clusters



Best Epsillon = 0.9
Best N_Samples = 50.0



Estimated Number of Clusters: 7.0

Best DBSCAN Scores for Cosine Distance Metric

Homogeneity score: 0.7896799513244787 Completeness score: 0.667782538810624 V_measure score: 0.7236336939202158

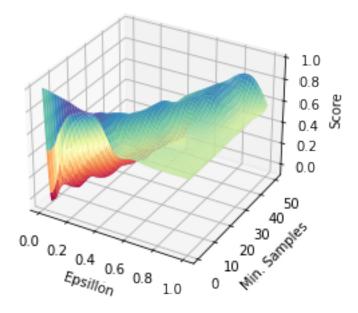
Adjusted Rand Index score: 0.7108732847743878

Adjusted Mutual Information Score: 0.7221232251101021

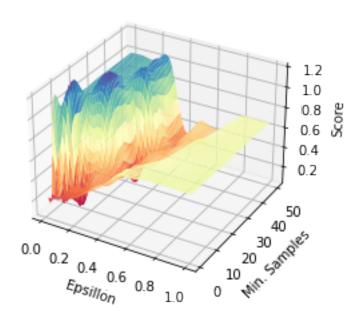
```
samples = [1,3,5,10,20,30,35,40,50]
metric_scores_DBSCAN_EUC = []
for eps in eps_vals:
   for sam in samples:
        DBSCAN_clust = DBSCAN(min_samples=sam, eps=eps)
        DBSCAN_clust.fit(bbc_umap)
        labels = DBSCAN_clust.labels_
        n_clusters_ = len(set(labels)) - (1 if -1 in labels else 0)
        metric_scores_DBSCAN_EUC.append([metrics.homogeneity_score(bbc.target,_
 →DBSCAN_clust.labels_),
                                     metrics.completeness_score(bbc.target,__
 →DBSCAN_clust.labels_),
                                     metrics.v_measure_score(bbc.target,__
 →DBSCAN_clust.labels_),
                                     metrics.adjusted_rand_score(bbc.target,__
 →DBSCAN_clust.labels_),
                                     metrics.adjusted_mutual_info_score(bbc.
 →target, DBSCAN_clust.labels_),
                                     n_clusters_,eps,sam])
# Visualize parameter effects on metrics
from mpl_toolkits.mplot3d import Axes3D
from scipy.interpolate import griddata
metricdata = np.array(metric_scores_DBSCAN_EUC)
metric_scores = ['Homogeneity', 'Completeness', 'V Measure', 'Adjusted Rand',
→'Adj. Mutual Info', 'Number of Clusters']
for i in range(0,6):
   plotdata = metricdata[:, [i,6,7]]
    z, x, y = zip(*plotdata)
    z = list(map(float, z))
    grid_x, grid_y = np.mgrid[min(x):max(x):100j, min(y):max(y):100j]
    grid_z = griddata((x, y), z, (grid_x, grid_y), method='cubic')
   fig = plt.figure()
    ax = fig.gca(projection='3d')
    ax.plot_surface(grid_x, grid_y, grid_z, cmap=plt.cm.Spectral)
    ax.set_xlabel('Epsillon')
    ax.set_ylabel('Min. Samples')
    ax.set_zlabel('Score')
    ax.set_title(str(metric_scores[i]))
```

```
plt.show()
# Find best parameter combination for epsillon and minimum samples
avs = []
for i in range(len(metricdata)):
→((metricdata[i][0]+metricdata[i][1]+metricdata[i][2]+metricdata[i][3]+metricdata[i][4])/
 →5)
    avs.append(a)
b = avs.index(max(avs))
print('Best Epsillon = ', metricdata[b][6])
print('Best N_Samples = ', metricdata[b][7])
cm = contingency_matrix(bbc.target, DBSCAN_clust.labels_)
rows, cols=linear_sum_assignment(cm, maximize=True)
plot_mat(cm[rows[:, np.newaxis], cols], xticklabels=cols, yticklabels=rows,_u
 \rightarrowsize=(5,5))
print("Estimated Number of Clusters: ", metricdata[b][5])
print()
print("Best DBSCAN Scores for Euclidean Distance Metric")
print("Homogeneity score: ", metricdata[b][0])
print("Completeness score: ", metricdata[b][1])
print("V_measure score: ", metricdata[b][2])
print("Adjusted Rand Index score: ", metricdata[b][3])
print("Adjusted Mutual Information Score: ", metricdata[b][4])
```

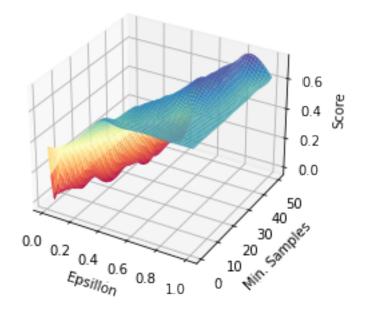
Homogeneity



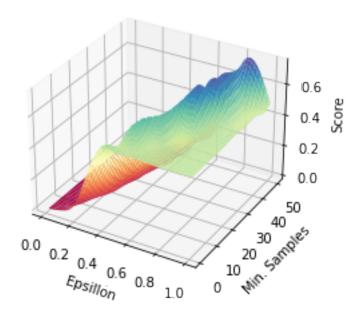
Completeness



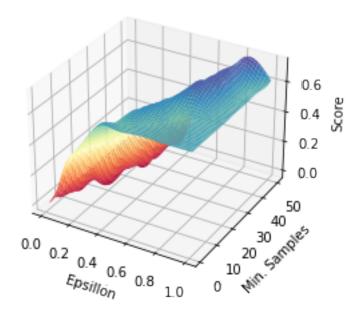
V Measure



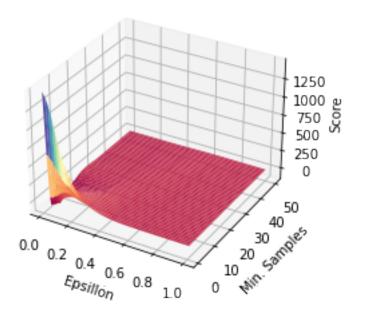
Adjusted Rand



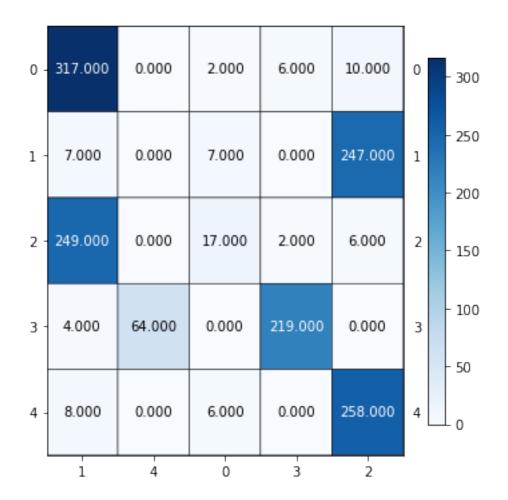
Adj. Mutual Info



Number of Clusters



Best Epsillon = 0.9
Best N_Samples = 50.0



Estimated Number of Clusters: 7.0

Best DBSCAN Scores for Euclidean Distance Metric

Homogeneity score: 0.7896799513244787 Completeness score: 0.667782538810624 V_measure score: 0.7236336939202158

Adjusted Rand Index score: 0.7108732847743878

Adjusted Mutual Information Score: 0.7221232251101021

[]: