PROJECT REPORT

The Implementation of DBSCAN and DBSCANRN with Tanimoto Similarity

"Data Mining"

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Introduction

Clustering is one of the most important tasks of both artificial intelligence and data mining. That is a way to group a set of data points in a way that similar data points are grouped together. Therefore, clustering algorithms look for similarities or dissimilarities among data points. Clustering is an unsupervised learning method so there is no label associated with data points. The algorithm tries to find the underlying structure of the data. There are different approaches and algorithms to perform clustering tasks which is of them is algorithms based on density clustering. Based on discussed in the lecture, we will implement algorithms of Density-Based Spatial Clustering of Applications with Noise (DBSCAN) and Density-based spatial clustering using reverse nearest neighbors (DBSCRN).

However, noted during project consultations, there are some inconsistencies in the paper presenting the pseudo-code of the DBSCRN algorithm [1]. For this reason, we will implement some modification of DBSCRN (which will be called DBSCANRN) instead of DBSCRN. In DBSCANRN, we will keep core point and non-core points definitions as given in the pseudo-code of DBSCRN. Namely, a point is considered as a DBSCRN (and by this DBSCANRN) non-core point if the number of their reverse k-nearest neighbors is less than k (lines 2-3 in the DBSCRN pseudo-code). Otherwise, it is considered as a core point (line 5 in the DBSCRN pseudo-code). We implement DBSCANRN in a way similar to NBC, except that instead of adding k+-nearest neighbors of a core point to a cluster, its reverse k-nearest neighbors will be added to the cluster.

Similarity Measures

The Tanimoto similarity is commonly used in bio-informatics or biology, and information retrieval to identify neighborhoods of sufficiently similar items or the k most similar things represented by real-valued vectors. The triangle inequality condition is frequently used to rapidly discover vectors that may belong to the targeted neighborhood of a given vector for metrics such as the Euclidean distance. However, the Tanimoto similarity and Tanimoto dissimilarity do not satisfy the triangle inequality property for real-valued vectors [2].

Considering this, another way to find a neighborhood with regard to Tanimoto similarity among real-valued vectors has been explained in [2]. We will adopt it, using bounds on vector lengths to calculate Tanimoto similarity ε -neighborhoods.

Property. Let u and v be non-zero vectors,
$$\alpha = \frac{1}{2} \left(\left(1 + \frac{1}{\varepsilon} \right) + \sqrt{\left(1 + \frac{1}{\varepsilon} \right)^2 - 4} \right)$$
, and $\varepsilon \in (0,1]$. If $T(u,v) \geq \varepsilon$, then $|v| \in \left[\frac{1}{\alpha} |u|, \alpha |u| \right]$. Their usefulness with an example illustrated in [2].

Datasets

We have 5 datasets:

- the toy dataset used in the class to present the execution of the NBC clustering algorithm (slide 51)
- 2. dim512 from http://cs.joensuu.fi/sipu/datasets/
- complex9 from https://github.com/deric/clusteringbenchmark/tree/master/src/main/resources/datasets/artificial
- 4. luto-t7-10k from https://github.com/deric/clusteringbenchmark/tree/master/src/main/resources/datasets/artificial
- letter from https://github.com/deric/clusteringbenchmark/tree/master/src/main/resources/datasets/real-world

Table 1. Used Datasets

No.	Dataset	Size	Dimension	Cardinality
0	lacture	12	2	<na></na>
1	dim512	1024	512	16
2	complex9	3031	2	9
3	cluto t7 10k	10000	2	10
4	letter	20000	16	26

measure labeled dataset based on size (num of observation), dimension (num of features), and cardinality (num of unique class). We also visualize the datasets in which datasets that have more than 2 dimensions will be transformed using principal component analysis (PCA) into 2 principal components, so that we can graph into 2-dimensional space. The graph of the datasets as follow:

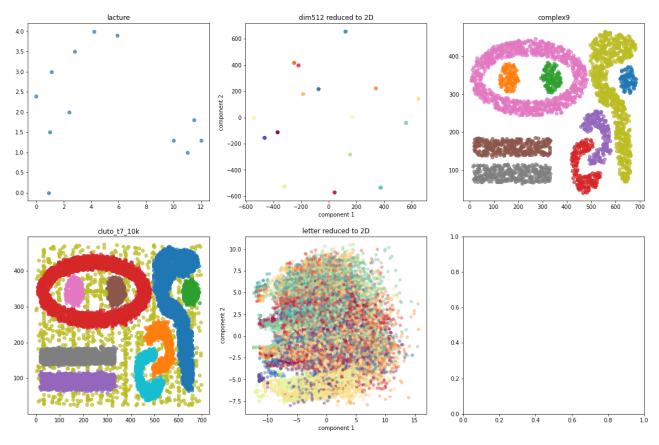


Figure 1. datasets

User's Manual

To run the program, open the Jupyter Notebook file named "DAMI_Project.ipynb", from this link: https://github.com/rauzansumara/dbscan-and-dbscanrn-tanimoto-similarity. This interactive Jupyter Notebook saved the results from previous run by us. Users can modify and run each cell from top to bottom to view the result of the code. Note that this action will replace previous output plots. The environment and library requirements as follow:

```
Requirements

python==3.8.5

numpy==1.21.4

pandas==1.3.5

sklearn==0.24.2

matplotlib==3.3.2
```

Experimental Results

We conducted our experiment using python programming language. Once you open the project folder, you will be presented with the following directory structure:

```
$ tree-project-folder
 datasets
   ├─ cluto-t7-10k.arff
    -- complex9.arff
   ├─ dim512.pa
    — dim512.ts
    — dim512.txt
     — lecture.csv
   letter.arff
 images
   ├─ datasets.png
   dbscan_cluto_t7_10k.png
   dbscan_complex9.png
   dbscan_dim512.png
   dbscan_lacture.png
   dbscan_letter.png
    dbscanrn_cluto_t7_10k.png
   dbscanrn_complex9.png
   ├─ dbscanrn_dim512.png
   optdbscanrn_letter.png
  - docs
   documentation [report]
  - save_file
   DEBUG_dbscanrn_cluto_t7_10k_D2_R10000_k10.csv
   DEBUG_dbscanrn_complex9_D2_R3031_k9.csv
   DEBUG_dbscanrn_lacture_D2_R12_k2.csv
   DEBUG_dbscanrn_letter_D16_R20000_k17.csv
   DEBUG_dbscanrn_cluto_t7_10k_D2_R10000_k10.csv
   <del>---</del> .....
   ├── STAT_dbscan.csv
    — STAT_dbscanrn.csv
   └── STAT_Opt_dbscanrn.csv
  - DAMI_project.ipynb
```

The *datasets* directory contains the data used. The *images* directory contains visualization graph of each experiment. We also have *docs* directory providing report file, and the most significant things are the *save_file* directory storing the DEBUG, OUTPUT, and STAT files. The results returned by a clustering algorithm for a given dataset and parameter values are saved in in 3 files:

<u>DEBUG files</u> - a file that contains important information for each point in the dataset on a separate line. Those are,

point id, minEps, maxEps, |Rk+NN|, identifiers if k+NN, identifiers of Rk+NN where:

- > maxEps Eps value calculated based on first k candidates for k+NN of the point
- > minEps the minimal value of Eps within which real k+NN of the point was found

<u>OUT files</u> - an output file that contains the following information on a separate line for each point in the dataset:

where:

- point id the position of the point in the dataset (before possible sorting),
- > x, y, ... dimension values
- TypeId either a core point (denoted by 1), or a border point (denoted by 0), or a noise point (denoted by -1)
- Clld a cluster identifier (which is a natural number) or -1 in the case of noise points.

<u>STAT files</u> - a file with the following statistics:

Dataset, Size, Dimension, Cardinality, Clld, Core_Point, Border_Point, Noise_Point,
Reading_Time, kNN/rkNN_Time, Clustering_Time, Save_Time, Total_Time, Purity_Index,
Rand_Index, DaviesBouldin_Index, Silhouette_Score

Implementation of DBSCAN

In this section, we did some set of experiments using DBSCAN algorithm as follow:

- ➤ lacture dataset DBSCAN(lacture, eps=0.80, MinPts=4)
- dim512 dataset DBSCAN(dim512, eps=0.80, MinPts=5)
- complex9 dataset DBSCAN(complex9, eps=0.9995, MinPts=5)
- cluto_t7_10k dataset DBSCAN(cluto_t7_10k, eps=0.99975, MinPts=6)
- ➤ letter dataset DBSCAN(letter, eps=0.99975, MinPts=7)

Once running the code, we obtained STAT file with following results:

Table 2. STAT file of the DBSCAN experiment

Droportios	Dataset					
Properties	lecture	dim512	complex9	cluto_t7_10k	letter	
Size	12	1024	3031	10000	20000	
Dimension	2	512	2	2	16	
Cardinality	<na></na>	16	9	10	26	
Clld	3	16	50	78	17	
Core_Point	5	1024	2160	7065	156	
Border_Point	3	0	306	793	0	
Noise_Point	4	0	565	2142	19844	
Reading_Time	0.004002	0.620927	0.036099	0.114168	0.452215	
Clustering_Time	0.001989	17.812995	51.045889	667.353465	2475.006246	
Save_Time	0.001999	0.075964	0.011995	0.032015	0.332090	
Total_Time	0.007991	18.509887	51.093983	667.499647	2475.790551	
Purity_Index	0.916667	1.000000	0.916529	0.933300	1.000000	
Rand_Index	0.893939	0.061584	0.565266	0.593961	0.999944	
DaviesBouldin_Index	1.905921	0.020462	1.255038	2.629512	1.117191	
Silhouette_Score	0.337355	0.986691	-0.089414	-0.307721	-0.332474	

As we can see from table above, we found 3 cluster in lacture dataset (include Noise). From 12 points, we got 4 core points, 3 border points, and 4 noise points. The total of running time is 0.007991 seconds. We obtained also the quality measurement such as Purity_Index, Rand_Index, DaviesBouldin_Index, and Silhouette_Score. We provided two graphs of each dataset to see the different between original clusters and predicted clusters. After running the code following results were obtained:

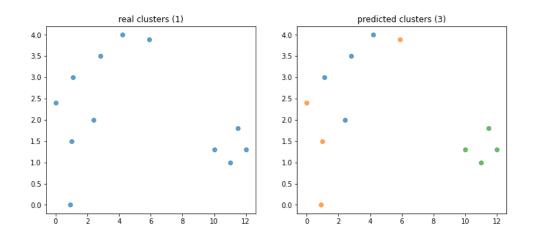


Figure 2. Visualization of Lecture Dataset (Left: Original and Right: Predicted Cluster)

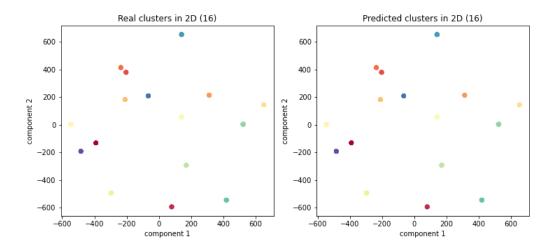


Figure 3. Visualization of dim512 Dataset (Left: Original and Right: Predicted Cluster)

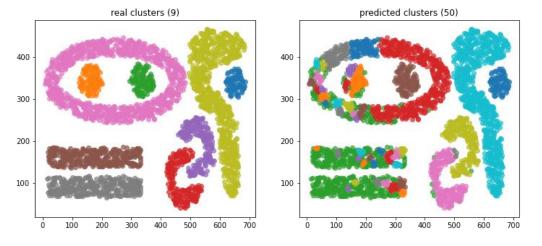


Figure 4. Visualization of complex9 Dataset (Left: Original and Right: Predicted Cluster)

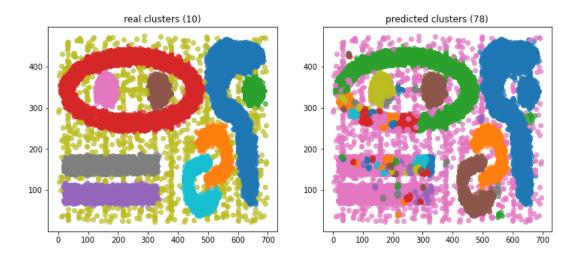


Figure 5. Visualization of cluto_t7_10k Dataset (Left: Original and Right: Predicted Cluster)

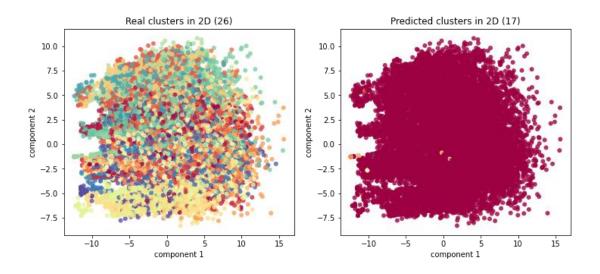


Figure 6. Visualization of letter Dataset (Left: Original and Right: Predicted Cluster)

Implementation of DBSCANRN

In Implementation of this algorithm both *non-optimized* and *optimized* version, we did some set of experiments as follow:

- lecture dataset DBSCANRN(lecture, k=2)
- dim512 dataset DBSCANRN(dim512, k=4)
- complex9 dataset DBSCANRN(complex9, k=9)
- cluto_t7_10k dataset DBSCANRN(cluto_t7_10k, k=10)
- letter dataset DBSCANRN(letter, k=17)

Once running the code, we obtained STAT file with following results:

Table 3. STAT file of the DBSCANRN experiment with Non-Optimized Version

Duonoution	Dataset				
Properties	lecture	dim512	complex9	cluto_t7_10k	letter
Size	12	1024	3031	10000	20000
Dimension	2	512	2	2	16
Cardinality	<na></na>	16	9	10	26
Clld	2	27	39	62	58
Core_Point	8	209	1753	5891	9896
Border_Point	4	815	1277	4077	10098
Noise_Point	0	0	1	32	6
Reading_Time	0.004002	0.619919	0.035098	0.113179	0.449201
kNN/rkNN_Time	0.002001	18.267762	55.351149	621.120478	2615.576711
Clustering_Time	0.000000	0.001000	0.004003	0.014990	0.044013
Save_Time	0.006001	0.087137	0.070135	0.253114	0.953630
Total_Time	0.012005	18.975818	55.460385	621.501762	2617.023555
Purity_Index	0.666667	0.795898	0.580666	0.597100	0.521900
Rand_Index	0.454545	0.344243	0.489176	0.488377	0.500591
DaviesBouldin_Index	0.307556	2.102505	0.815071	1.561321	1.225597
Silhouette_Score	0.713755	0.708637	0.210441	-0.118104	-0.207850

Table 4. STAT file of the DBSCANRN experiment with Optimized Version

Droportion	Datasets					
Properties	lecture	dim512	complex9	cluto_t7_10k	letter	
Size	12	1024	3031	10000	20000	
Dimension	2	512	2	2	16	
Cardinality	<na></na>	16	9	10	26	
Clld	2	27	39	62	2	
Core_Point	8	209	1753	5891	11614	
Border_Point	4	815	1277	4077	8386	
Noise_Point	0	0	1	32	0	
Reading_Time	0.009998	0.620907	0.035098	0.113179	0.449209	
kNN/rkNN_Time	0.000999	5.269698	7.095079	31.629475	1664.005344	
Clustering_Time	0.000000	0.000998	0.002089	0.018149	0.053006	
Save_Time	0.014326	0.088201	0.066320	0.316230	0.938037	
Total_Time	0.025323	5.979804	7.198586	32.077034	1665.445597	
Purity_Index	0.666667	0.795898	0.580666	0.597100	0.580700	
Rand_Index	0.454545	0.344243	0.489176	0.488377	0.512648	
DaviesBouldin_Index	0.307556	2.102505	0.815071	1.561321	0.770860	
Silhouette_Score	0.713755	0.708637	0.210441	-0.118104	0.199443	

As we can see from table 3 (non-optimized version) and table 4 (optimized version), we found the same number predicted clusters and type of point (core, border or noise) on 4 datasets, such as lecture, dim512, complex9, and cluto_t7_10k datasets between non-optimized and optimized version.

Unfortunately, we do not know why it showed different result on letter dataset between non-optimized version and optimized version. But apart from it, the comparison of DBSCANRN non-optimized and optimized version proven that the optimized version can reduce computation time up to 19 times faster than original non-optimized version. Meanwhile, the computation time does not really much different when we apply in small dataset like lecture dataset. We provided two result graphs of each version to see the different between non-optimized and optimized version. After running the code following results were obtained:

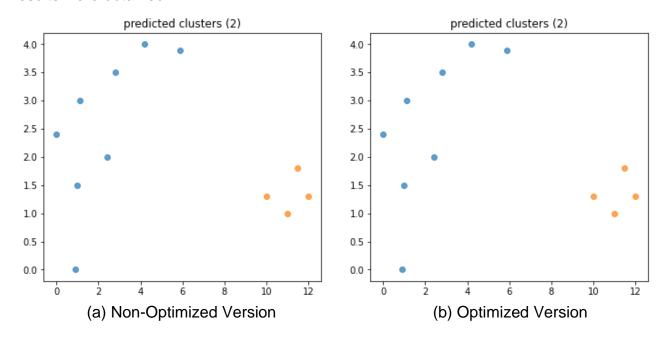


Figure 7. Visualization of Lecture Dataset

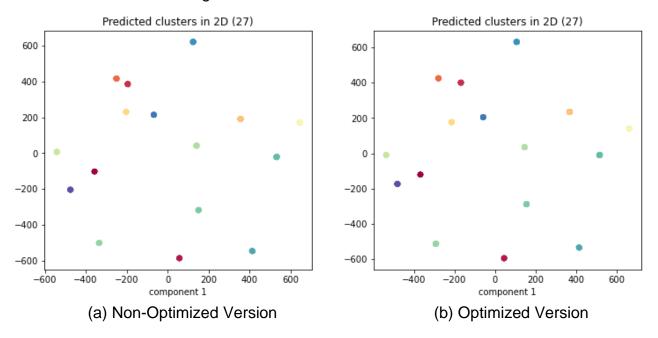


Figure 8. Visualization of dim512 Dataset

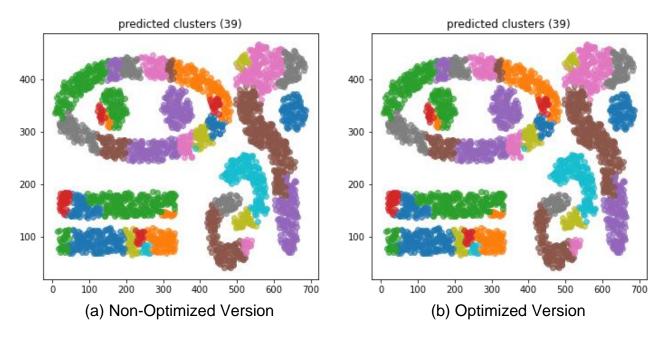


Figure 9. Visualization of complex9 Dataset

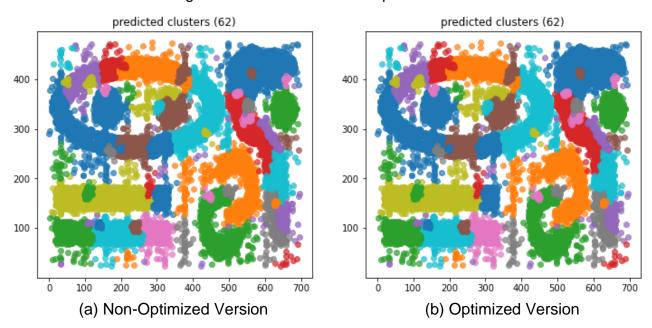


Figure 10. Visualization of cluto_t7_10k Dataset

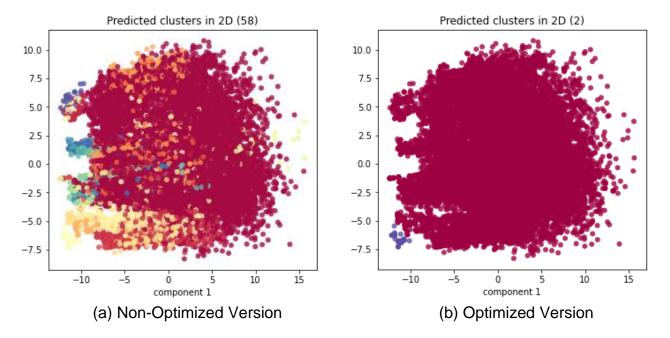


Figure 10. Visualization of letter Dataset

Conclusion

- in DBSCAN, computation time is affected by epsilon radius and size of dataset,
- in DBSCAN, larger *epsilon* creates more members of a cluster, while larger *minPts* usually creates a smaller number of clusters,
- when running program/code with the same parameters on DBSCANRN both versions (non-optimized and optimized version), our implementation of optimized one successfully reduced computation time up to 19 times faster compared to non-optimized version, and
- We obtained different number of found clusters only in letter dataset when comparing non-optimized and optimized version of DBSCANRN. It could be caused of code or characteristics of the data itself; we need to investigate in the future.

References

- [1] Chowdhury, S. and Amorim R.C., (2019) An efficient density-based clustering algorithm using reverse nearest neighbour. The Computing Conference 2019: London
- [2] Kryszkiewicz M. (2021) Determining Tanimoto Similarity Neighborhoods of Real Valued Vectors by Means of the Triangle Inequality and Bounds on Lengths.
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- [4] Kryszkiewicz M. (2014) Using Non-Zero Dimensions and Lengths of Vectors for the Tanimoto Similarity Search among Real Valued Vectors. Springer International Publishing Switzerland 2014