Arbitrary Shape Cluster Summarization with Gaussian Mixture Model

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Abstract: One of the main concerns in the area of arbitrary shape clustering is how to summarize clusters. An accurate representation of clusters with arbitrary shapes is to characterize a cluster with all its members. However, this approach is neither practical nor efficient. In many applications such as stream data mining, preserving all samples for a long period of time in presence of thousands of incoming samples is not practical. Moreover, in the absence of labelled data, clusters are representative of each class, and in case of arbitrary shape clusters, finding the closest cluster to a new incoming sample using all objects of clusters is not accurate and efficient. In this paper, we present a new algorithm to summarize arbitrary shape clusters. Our proposed method, called SGMM, summarizes a cluster using a set of objects as core objects, then represents each cluster with corresponding Gaussian Mixture Model (GMM). Using GMM, the closest cluster to the new test sample is identified with low computational cost. We compared the proposed method with ABACUS, a well-known algorithm, in terms of time, space and accuracy for both categorization and summarization purposes. The experimental results confirm that the proposed method outperforms ABACUS on various datasets including syntactic and real datasets.

1 INTRODUCTION

Nowadays, a large volume of data is being generated which is even difficult to be captured and labelled. A large volume of data is mainly generated in stream and real time applications, in which, data is generated rapidly, and cannot be stored bit by bit. As a result, analysing and labelling such kind of data is a main challenge (Guha et al., 2003)(Bifet et al., 2009) (Charu et al., 2003). Lack of labelled data draws attention to the application of clustering to produce labelled data. The choice of clustering algorithm strongly depends on data characteristic. In small and noise free environment classical clustering method like k-mean and k-median are commonly used. However, in most of applications, there is no knowledge about the number of clusters while the shapes of clusters are non-convex and arbitrary. In this case, density-based and grid-based clustering methods are used. Using a clustering method that generates arbitrary shape clusters is theoretically ideal but the representation and analysis of each cluster still causes many problems. For full representation of arbitrary shape clusters, all the samples of clusters should be preserved which is impractical in many applications. In case of using clustering in online applications, each cluster can be representative of a specific pattern. These patterns need to be kept for a long time and keeping the full representation of the complex patterns tend to be impractical. In this case, summarization and extracting the key features of clusters are necessary.

Another application of clustering is categorizing the unlabelled data. In this case, a set of clusters is created and each cluster receives a label according to its own samples. Each new sample is compared to the clusters and the closest one is chosen as a cluster that the new sample belongs to. In case of k-means, the distance of a new sample to the centre of the cluster is calculated, then, if the distance is less than the radius of a cluster, new sample is attached to that cluster. K-means and partition-based clustering methods are sensitive to noise and cannot detect clusters with arbitrary shape. This approach has been

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studied in different applications but most of the attention was towards non-arbitrary shape clustering like k-means (Mohammadi et al., 2014)(Gaddam et al., 2007). Arbitrary shape clustering methods preserve all samples for each cluster and to find the closet cluster to the new sample, distance of new samples to all cluster members is calculated. It is obvious that using such an approach is time consuming. The other way to find the closet cluster is to create a boundary for each cluster. If the new sample is inside the boundary of a cluster then the new sample belongs to that cluster. Finding the boundary of arbitrary shape clusters, especially in high dimensional problems, is a complex and time consuming process. Moreover, it is necessary to save too many faces to just keep borders of cluster created by convex in higher dimension, which grows exponentially with dimension (Kersting et al., 2010)(Hershberger, 2009).

In this paper, we propose a new approach that fulfils the mentioned requirements. We propose a summarization approach to summarize arbitrary shape clusters using Gaussian Mixture Model (GMM). In our approach, we first find the core objects of clusters and then we consider these core objects as centres of GMM and represent a cluster with a GMM. Since, GMM-based method keep all statistical information of each cluster, it summarizes each cluster in a way that we can use it for pattern extraction, pattern matching, and pattern merging. Moreover, this model is able to classify new objects. Using GMM, each new test sample is fed into the GMM of a cluster, and if the membership probability to a cluster is more than a threshold, the object is attached to that cluster.

The structure of the paper is as follows: In Section 2, we review related work on arbitrary shape clusters and summarization approaches. In Section 3, we explain the general structure of the proposed algorithm for summarization. In Section 4, we present some discussions about the features of the proposed method. In Section 5, we explain the complexity of algorithm in more detail. Section 6 presents the experimental results of the proposed algorithm in comparison with well-known summarization algorithms. Finally, the conclusion and future work are presented in Section 7.

2 RELATED WORK

There are various algorithms available for clustering, which are categorized into four groups; partition-based, hierarchical, density-based and spectral-based clustering (Han, 2006). K-means is one of the famous algorithms in the area of partition-based clustering. However, using a centre and radius makes the shape of clusters spherical which is undesirable in many applications. In hierarchical clustering methods such as Chameleon data is clustered in hierarchical form but still with spherical shape that is undesirable. Moreover, tuning the parameters for methods like Chameleon is still difficult (Karypis et al., 1999). Spectral clustering; STING (Wang et al., 1997) and CLUIQE (Agrawal et al. 1998) are able to create arbitrary shape clusters but the major drawback of these methods is the complexity of creating an efficient grid. The size of grid varies for different dimensions and setting different grid sizes and merging the grids to find clusters are difficult. These difficulties make the algorithm inaccurate in many cases. In the area of arbitrary shape clustering, density-based methods are more interesting and DBSCAN (Ester et al., 1996) and DENCLUE (Hinneburg et al., 1998) are the most famous ones. In density-based methods, clusters are created using the concept of connecting dense regions to find arbitrary shape clusters. Based on prevalence of real time applications, there is more interest to make these algorithms fast for streaming applications (Guha et al., 2003)(Bifet et al., 2009)(Charu et al., 2003).

Summarization is the solution to ease the complexity of arbitrary shape clustering methods. The naïve way to represent an arbitrary shape cluster is to represent each cluster with all cluster members. Obviously, this approach is neither practical nor does it reflect the cluster properties. In k-means a simple representation using a centre and radius summarize the cluster. It is clear that this summarization does not capture how data is distributed in the cluster.

There are different ways to summarize arbitrary shape clusters (Yang et al., 2011)(Cao et al., 2006)(Chaoji et al., 2011). These algorithms use the general idea behind the clustering methods for arbitrary shape clusters. In the area of summarization, the idea is to detect dense regions and summarize the regions using core objects. Then, a set of proper features is considered to summarize the dense regions and their connectivity. In (Yang et al., 2011) a grid is created for each cluster and based on the idea of connecting dense regions, the core or dense cells with their connections and their related features are kept. In all summarization approaches, these features play crucial role. In (Yang et al., 2011) location and range of values and status connection vector are kept however, it has some

draw backs. First, creating grids on each cluster is time consuming. Second, considering all grids needs spending a lot of time and space which is impractical in many cases. Cao et al. (Cao et al., 2006) use the idea of finding core objects to generate the cluster summery. The most significant drawback of this work is that the number of core objects is large and in some cases, it is equal to the number of input samples. Moreover, a fixed radius specifies the neighbourhood that does not represent the distribution of objects in each cluster (Cao et al., 2006). Chaoji et al. represent a density-based clustering algorithm named ABACUS for creating arbitrary shape clusters (Chaoji et al., 2011). The summarization part in their approach is based on finding the core objects and the relative variance around the objects. In most of arbitrary shape clustering methods, we need two parameters; number of neighbours and a radius. The most interesting and noticeable part in Chaoji et al. work is that the number of neighbours is the only parameter for their algorithm and they generate radius using data distribution. The significant drawback in their work is that the algorithm may generate many core points.

In all above mentioned summarization approach, focus is on preserving the cluster members and they don't consider any usage of clustering for classification purpose. Clustering approaches are created to summarize data and categorized the input data into some groups, but it can also be used as a pre-processing phase of classification task (Ester, 1996)(Hinneburg, 1998). Each cluster has a label and for each new object the closest cluster is found and the object gets the label of that cluster. The mentioned approaches like k-means summarize the clusters, but they still consider the concept of circles around the core objects to find the closest cluster which is inaccurate. Graph-based structures are more accurate but they are time consuming. Anomaly detection is one of the applications of clustering in classification (HE, 2003)(Borah and Bhattacharyya, 2008)(Gaddam, 2007)(Mohammadi, 2014). In this area, those objects which are outside of cluster boundaries are considered as anomaly. In previous works, clusters are generated using k-means and summarized with a centre and a radius which is inaccurate.

In this paper, we present an approach to summarize clusters using Gaussian Mixture Models. Our approach covers both areas; it is a good representative of a cluster and it can be used for classification purpose.

3 GENERAL STRUCTURE OF THE SGMM ALGORITHM

The main idea behind the density-based clustering algorithms is to connect dense regions to create a cluster. With this idea, a set of core objects are detected and they are connected to each other using the shared neighbourhood.

DEFINITION 3.1. *In dataset D for a given k and radius r, an object o_i is a core object if*

$$\{\forall x \in D, o_j \in C | \, \|\{d(o_i, x) < r\}\| \geq k\}$$

Where *C* is set of core objects and (|| ||) shows the number of objects with distance less than *r* from the core object. Detection of *k* and *r* is critical in identifying good clusters.

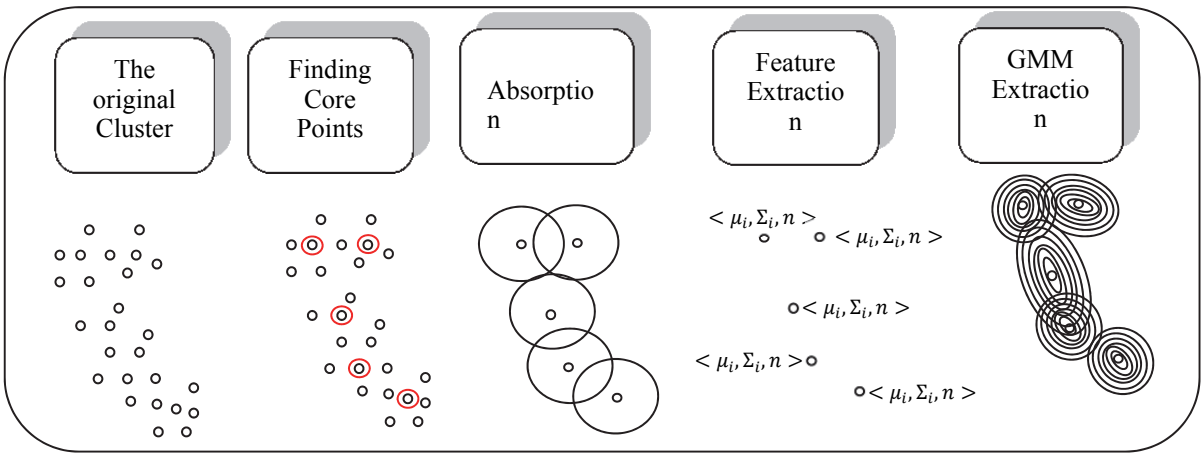

Figure 1: Structure of the SGMM method for cluster summarization.

The concept of core objects has been used to generate the backbone of a cluster and generate summery of the cluster. In some recent works (Chaoji, 2011) using a given *k* and core objects idea, they detect the backbone of cluster. In this paper, we use the same idea but with more concentration on decreasing the number of core objects and the time of detection of these core objects. The general idea for summarization of each cluster in this paper is based on Gaussian Mixture Model. Since Gaussian Mixture Model preserves the distribution of data using a set of Normal Distributions, it is a good candidate for summarizing any cluster. Therefore, we can apply simply EM algorithm on each cluster to find its GMM representative, but the main concern is to find the number of GMM components. In our proposed approach, we first find the number of GMM components, and then we find appropriate GMM for each cluster. Our method, Summarization Based on Gaussian Mixture Model (SGMM), has four main steps that are depicted in Figure 1. First, a set of objects called core objects are detected. These objects are representative of a cluster and can generate the original cluster as needed. After detection of backbone objects, there comes the absorption step, where the objects attached to the core object are absorbed and represented by core objects. Then by introducing a new feature set for each object, cluster is summarized while its original distribution is preserved. Finally, each cluster is presented as a GMM. In the following sections, we describe each step in more detail.

3.1 Finding Core Objects

In this step, we find the core objects that create the backbone of each cluster. There are different ways to find the core objects in each cluster, but time is the main concern in this case. In this step, we consider a radius based on which we find the neighbours for all objects for each cluster. Then we sort all objects based on the number of neighbours. The first core object is the one with the maximum number of neighbours. All the neighbours of this object are removed from the list of possible core objects. With this approach, we reduce the overlap of core objects as much as possible. Among the remaining objects, we find the next objects with maximum number of neighbours. We label this object as a core object and remove neighbours of this object from further consideration. In this way with a heuristic approach, we find the core objects located in dense regions and all parts of cluster are covered using them. The new definition of the core objects with our methodology is presented as follow.

DEFINITION 3.1.1. *(Core Object) A core object is the object that has the maximum number of neighbours in comparison with other objects in its neighbourhood. The core object is not in the neighbourhood of another core object.*

$$\{\forall c_i \in C | i \neq j , d(c_i, c_j) > r\}$$

Where *i* and *j* are the index for different core points and *d* refers to the distance of two objects.

3.2 Absorption and Cluster Feature Extraction

The goal of summarization is to find a good representative of each cluster and core objects are the only objects that we preserve in each cluster while the rest of the objects in the cluster are removed. After finding all core objects in each cluster, the next step is to define a cluster using core objects. It is obvious that considering only core objects cannot be a good representative of each cluster. The core objects have to be accompanied with a set of features related to the cluster to represent a cluster. This is why we define a set of features for each core object which are good representative for distribution around each core object.

DEFINITION 3.2.1: *(Core Object Feature) (CF) Each core object is represented by a triple* $\langle c_i, \Sigma_i, \omega_i \rangle$.

In this definition c_i is the core object and Σ_i is the covariance calculated using the core object and all objects in its neighbourhood. $\omega_i = n/CS$ is the weight of core object, *n* is the number of objects in the neighbourhood of core object c_i , and *CS* is cluster size. ω_i shows the proportion of objects which are in the neighbourhood of the core object. Using the features for each core object, we estimate samples scattering around each core object without keeping the entire samples in the neighbourhood.

3.3 GMM Representation of Clusters

After finding the core objects and all necessary features for each cluster, we generate a Gaussian Mixture Model for each cluster.

DEFINITION 3.3.1. *A Gaussian Mixture Model is a combination of a set of normal distributions. Given*

feature space $f \subset R^d$, a Gaussian Mixture Model $g: f \to R$ with n component is defined as:

$$g(x) = \sum_{i=1}^{n} w_i \mathrm{N}_{\mu_i \Sigma_i}(x)$$

$$\mathrm{N}_{\mu_i \Sigma_i}(x) = \frac{1}{\sqrt{(2\pi)^d |\Sigma_i|}} e^{-\frac{1}{2}(x-\mu_i)\Sigma_i^{-1}(x-\mu_i)^T} \quad (1)$$

Based on all $CF_i, i = 1 \cdots m$, a GMM is defined over a cluster. Each component for GMM is created using a core object, its covariance and the weights. In the formula in equation (1), μ_i is the centre of the i^{th} GMM component, which is set to the coordination of core object and therefore, $\mu_i = c_i$. The covariance is set to covariance of i^{th} core object covariance, that is Σ_i. The weight for each component is the weight of core object and as a result w_i in the equation (1) is set to the weight of core object c_i which is ω_i.

The main goal of cluster summarization is to present a cluster in a way that it keeps the overall distribution of the cluster while reducing the number of objects in the cluster which is critical in stream data mining and online environment. To use a cluster as a class in classification, we need to find the closest cluster to a new incoming sample. In an arbitrary shape cluster, all objects represent a cluster and to find a right cluster for a new sample, it has to be compared with all objects in a cluster. GMM-based summarization fulfils both requirements for both applications; summarize data for the real time applications; generate a proper representation for each cluster to use them in classification approach.

4 GAUSSIAN MIXTURE SUMMARIZATION

As each GMM consists of a set of Gaussian distributions, finding the number of GMM components is a challenging issue. In this paper, we proposed a solution to calculate number of components for each cluster by finding the core points. In our approach, number of components for GMM is set to the number of core objects in a cluster.

There are two essential features that should be considered. Each summarization technique has to preserve the original shape and the distribution of the data. Summarization of data using GMM has both characteristics. In SGMM some objects are selected in a way that they follow the general structure of data. Not only finding the core object is easy and fast but also it still follows the shape of the cluster. The algorithm starts from dense regions in the cluster and then goes to the most scattered part with consideration of covering all data objects in the cluster. In the SGMM method, the core objects are the ones which are in the center of dense regions and they cover all data. Therefore, for each region a representative object is chosen and then collection of these representative objects presents the entire cluster. Figure 2, shows the core objects generated for a cluster. This figure shows that the generated core objects follow the general structure of the cluster.

Figure 2: Cluster with the core objects.

Finding the general structure of a cluster using a set of objects is not enough for achieving high accuracy in classification applications. In many applications, we would like to keep the original or summery of data for more detail investigation. Saving the core objects cannot reveal any information about the distribution of data. Therefore, we need to find distribution information around the core objects and as we mentioned above, we capture such information using the covariance around core objects. Using such information, we are able to regenerate the original data.

Finally, we need to know the contribution of each core object in generation of data. We find out the weight of each core object based on the number of neighbours that each object has. With knowing the location of core objects, related covariance and the weights, we summarize data using a GMM with the minimum loss of information. The relative GMM for a cluster is depicted using contour plots in Figure 2.

5 COMPUTATIONAL COMPLEXITY

In stream data mining application, using a fast algorithm is a critical requirement. In this section, we examine the computational complexity of the proposed algorithm.

In the first step of algorithm, we find the neighbours of each object; with *N* as the number of objects it takes $O(N^2)$ to find neighbours. Then, we need to sort the objects and absorb neighbours of the core objects that takes *O(NlogN)*. In absorbing step, we find the number of neighbours for each core object and its related variance. We consider *O(NlogN)* for the second part of the proposed method which is the worst case.

So, the final complexity of the proposed algorithm is $O(N^2 + NlogN)$. In the grid-based method a considerable amount of time is spent to create the grid and further investigation to find core grids and connecting them. In some other density-based methods such as ABACUS there are iterations to find core objects and to reallocate them which are time consuming. The output of SGMM is almost the same as that of ABACUS but the proposed method is not time consuming in comparison to ABACUS.

The experiment results show that our heuristic leads to the same result with one step. In the next section, we develop a comparison between the required time of our method and the one for ABACUS. The results illustrate that the proposed method outperforms ABACUS in term of time.

6 EXPERIMENTAL RESULTS

We use both synthetic data set and the real data set in our experiments. Time and space complexity and the goodness of clustering are three different criteria that we evaluate our algorithm based on. Moreover, we set up different experiments to see the efficiency of our algorithm for categorizing new samples. We compared our algorithms with ABACUS (Chaoji, 2011) which is one of the well-known summarization methods in literatures. All experimental results are generated in Matlab running on a machine with Intel CPU 3.4 GHz and 4GB memory.

The first data set is a synthetic dataset as shown in Figure (3). We use this dataset to visualize the effectiveness of the proposed algorithm.

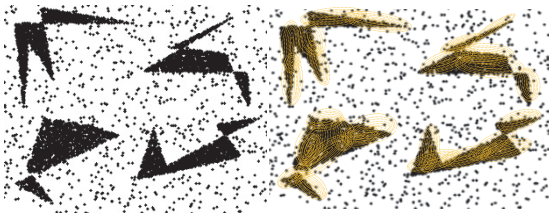

Figure 3: Synthetic data with related GMMs.

This figure shows 4 clusters and set of objects used in testing phase. In this Figure we see each cluster is represented by a GMM depicted with a contour plot. KDD dataset and some other UCI datasets are considered in our experiments to evaluate the accuracy of our algorithm on some real datasets.

6.1 Clustering Goodness

To show the efficiency of SGMM method, we set up experiments in which we summarize the dataset using core objects then we used these core objects and their related variance and weights to regenerate the dataset. The difference between the first dataset and regenerated dataset by core objects shows the strength of summarization algorithm. Experimental results show that SGMM method summarize dataset better than ABACUS method. To visualize the results, we generated a synthetic data with four clusters. Figure 4 from left to right shows the original dataset, core objects of each cluster and the dataset set regenerated using core objects using ABACUS and SGMM method. This figure shows that the core objects follow the original structure of the clusters and the regenerated clusters are similar to the original ones. This figure depicts the ability of SGMM method, and it shows that the summery generated by SGMM regenerates the original dataset better in comparison to ABACUS method. The difference between the accuracy of summarization of ABACUS and SGMM is clearer in the cluster with star shape. SGMM summery regenerate the data with star shape, but ABACUS method could not regenerate the same shape. These figures are just for visualization the result and to show the efficiency of the algorithm, we use Dunn and DB index.

The Dunn index (Dunn, 1979) is a validity index which identifies compact and well-separated classes defined by equation (2) for a specific number of classes:

$$D_{nc} = \min_{i=1,\dots,nc} \left\{ \min_{j=i+1,\dots,nc} \left(\frac{dist(c_i, c_j)}{\max_{k=1,\dots,nc} diam(c_k)} \right) \right\} \quad (2)$$

Where *nc* is the number of classes, and $dist(c_i,c_j)$ is the dissimilarity function between two classes C_i and C_j. The large values of the index indicate the presence of compact and well-separated classes. In our experiments, we first calculate the Dunn index for the original datasets. Then, summarize the dataset and re-generate the data again using the final core objects of GMMs and we find out the Dunn index for regenerated data. Finally, we get the difference between Dunn index of original dataset

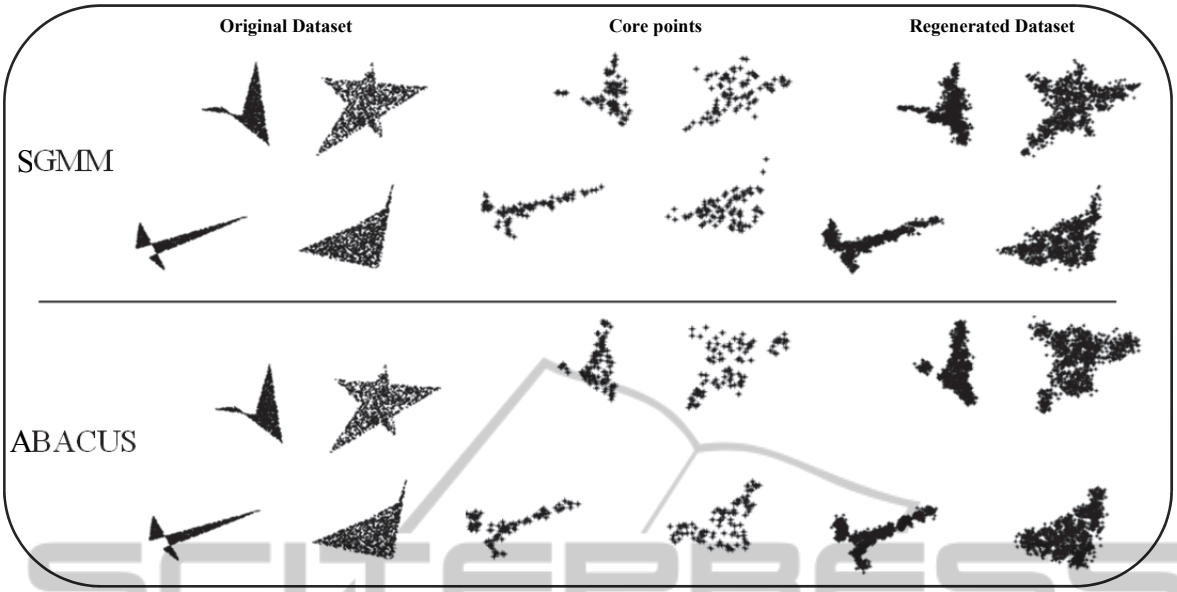

Figure 4: The first dataset is the original dataset, the second one is the core points and the third one is regenerated dataset using the core points based on ABACUS and SGMM methods.

and the one which is regenerated. Using this experiment, we want to evaluate the regeneration ability of the SGMM method. In other words, we want to show how the SGMM method regenerates the data which follows the shape and distribution of the original data. Table 1 shows the results of this experiment. Each value in this table is the average of 30 independent runs.

Table 1: Difference of Dunn Index for original and regenerated dataset.

Dataset\Index	SGMM	ABACUS
Synthetic Data	0.02358	0.076932
KDD	0.027701	0.034321
Segment	0.008841	0.014636

The closer the value to zero, the better the result we get. Based on the difference of Dunn index table, SGMM always outperforms ABACUS by far. For example in case of Segment dataset, the difference between the dataset generated by SGMM and the original dataset is almost zero while this value for ABCUS is around 0.01. This experiment illustrated that SGMM follow the sample distribution of the original dataset better than ABACUS. The next step is to repeat the experiment using DB cluster index.

The second measurement we used is DB index (Davies and Bouldin, 1979) which is a function of the ratio of the sum of within cluster scatter to between-cluster separation. DB index is defined as in equation (3).

$$DB = \frac{1}{n}\sum_{i=1}^{n} \max_{i \neq j}\left[\frac{S_n(Q_i)+S_n(Q_j)}{S_n(Q_i,Q_j)}\right] \quad (3)$$

n is the number of clusters, S_n is the average distance of all objects of the cluster to their cluster centre, and $S_n(Q_i,Q_j)$ is the distance between clusters centres. Hence, the ratio is small if the clusters are compact and far from each other. Consequently, Davies-Bouldin index will have a small value for a good clustering. As what we did for Dunn index, we find the distance of the original and regenerated data. Table 2 shows the experimental results of DB index on different datasets.

Table 2: Difference of DB Index for original and regenerated dataset.

Dataset\Index	SGMM	ABACUS
Synthetic Data	0.006638	0.017609
KDD	0.195763	0.232498
Segment	0.0429764	0.0546819

Result based on Dunn and DB index shows that the SGMM method generates more accurate summery of dataset and in case of regenerating the original dataset SGMM follows the original pattern better.

6.2 Clustering Accuracy

As mentioned above, clustering is used in classification as a preprocessing step. The application we focused on, in this part, is anomaly detection. We first cluster the normal data into some clusters then we consider these clusters as normal behaviours. The membership probability for each new sample is calculated. If the sample belongs to a cluster, it is normal otherwise it is an attack. For this purpose, we consider KDD dataset and we find out the accuracy of attack detection. Using GMM in classification is our proposed method and to the knowledge of author there is no other work similar to our approach. The only approach that has been used in an arbitrary shape cluster to label new sample is to consider all member of clusters which is the Naïve approach. In our experiment, we created GMM using our set of core objects and ABACUS core objects. As we mentioned, we created the core objects using a simple method that spends less time and space. Therefore, there can be a doubt about the accuracy of categorizing using the clusters using our method? The experiment shows that our approach does not decrease the accuracy.

The ROC curve based on detection rate and false alarm rate for KDD dataset depicted in Figure 5. It shows while the ABACUS is more time consuming and find out too many core objects our method has the same and sometimes better accuracy. Figure 6 shows the result for synthetic dataset. It shows that the accuracy of our approach is better than Naïve and ABACUS method. The result for synthetic dataset set shows that in spite of using clustering in categorizing new objects the accuracy is still good and comparable to the accuracy of classification methods.

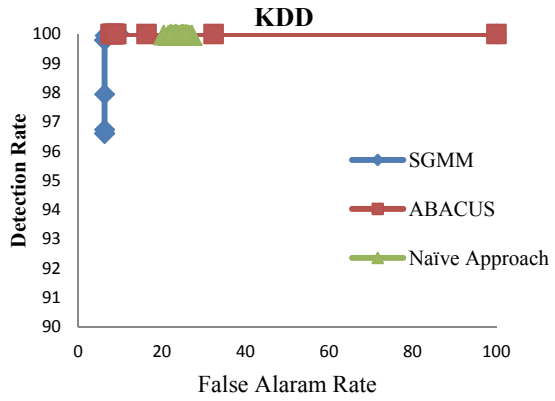

Figure 5: ROC curve for KDD dataset.

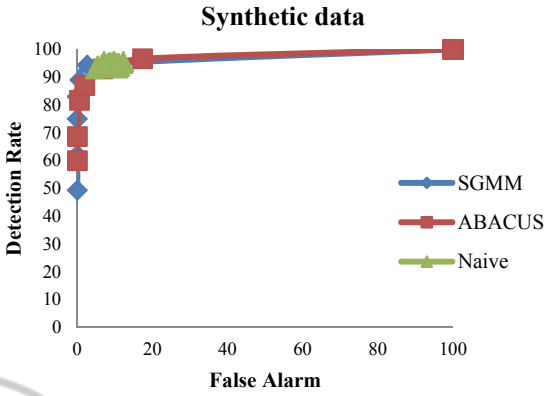

Figure 6: ROC curve for Synthetic dataset.

Table 3 shows more results on some other datasets, which confirms that summarizing and clustering data using SGMM method outperforms ABACUS method. The first value in Table 4 is the false alarm rate and the second one is the detection rate.

Table 3: Accuracy of SGMM and ABACUS method in anomaly detection.

Dataset\Index	SGMM	ABACUS
Synthetic Data	(2,94)	(1.9,87)
KDD	(6.301, 100)	(9.041, 100)
Segment	(33.83,94.39)	(32.82,94.54)

6.3 Algorithm Complexity

Two main concerns in the area of summarization are the time spent to find the summery of a cluster and the number of objects preserved for it. Table (4) shows the space and the time spent to find the summery of a cluster for different datasets. This table shows that SGMM decreases time and space to summarize each cluster. It is clear that SGMM method uses only one iteration to find core objects while ABACUS method runs many iterations to find out the core objects and therefore, SGMM is faster. On the other side, SGMM summarizes cluster with less number of core objects. The reason lies in the SGMM algorithm in finding core objects. If an object is core object, all objects in its neighborhood are removed from possible set of core objects and we do not consider them. In this case, we are able to decrease the time and space complexity both. Table 4 and 5 shows the time and space used by SGMM and ABACUS and it shows that using SGMM we decrease time and space considerably.

Table 4: Number of core objects for ABACUS and SGMM methods.

Dataset\Space complexity	SGMM	ABACUS
Synthetic Data	51	120
KDD	46	272
Segment	52	55

Table 5: The time complexity for ABACUS and SGMM methods.

Dataset\Time complexity	SGMM	ABACUS
Synthetic Data	388 sec	1261 sec
KDD	222 sec	836 sec
Segment	7 sec	19 sec

7 CONCLUSIONS

In this paper, we presented a new approach for summarizing the arbitrary shape clusters. Our proposed algorithm, named SGMM, represents each cluster by a Gaussian Mixture Model (GMM) using sets of core objects. Each GMM is a representative of the distribution of whole members in a cluster. Moreover, SGMM is able to find the closest cluster for each new incoming sample. The experimental results based on Dunn and DB index confirms that the distribution of clusters is preserved after summarization. Additionally, in case of classification, the accuracy using SGMM is acceptable. Our proposed algorithm also exhibits a low computational cost which makes it a suitable approach for clustering stream data. Extending the density based clustering to make it faster is a part of our future work.

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