

# A Modification on K-Nearest Neighbor Classifier

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**Abstract**-K-Nearest Neighbor (KNN) classification is one of the most fundamental and simple classification methods. When there is little or no prior knowledge about the distribution of the data, the KNN method should be one of the first choices for classification. In this paper a modification is taken to improve the performance of KNN. The main idea is to use robust neighbors in training data. This modified KNN is better from traditional KNN in both terms: robustness and performance. The proposed KNN classification is called Modified K-Nearest Neighbor (MKNN). Inspired from the traditional KNN algorithm, the main idea is to classify an input query according to the most frequent tag in set of neighbor tags. MKNN can be considered a kind of weighted KNN so that the query label is approximated by weighting the neighbors of the query. The procedure computes the frequencies of the same labeled neighbors to the total number of neighbors. The proposed method is evaluated on a variety of several standard UCI data sets. Experiments show the excellent improvement in the performance of KNN method.

**Keywords**- MKNN, KNN Classification, Modified K-Nearest Neighbor, Weighted K-Nearest Neighbor, Neighbor Validation.

## I. INTRODUCTION

Nowadays, recognition system is used in many applications which are related to different fields that have different nature. Pattern recognition is about assigning labels to objects which are described by a set of values named attributes or features. Current research builds upon foundations laid out in the 1960s and 1970s [1].

There are three major types of pattern recognition trends: unsupervised, semi-supervised and supervised learning. In the supervised category, also called classification or regression, each object of the data comes with a pre-assigned class label. In other hand, there is a teacher saying the true answer. The task is to train a classifier to perform the labeling, using the teacher. A procedure which tries to leverage the teacher's answer to generalize the problem and obtain his knowledge is learning algorithm. Most often this procedure cannot be described in a human understandable form, like Artificial Neural Networks classifiers. In these cases, the data and the teacher's labelings are supplied to the machine to run the procedure of learning over the data. Although the classification knowledge learned by the machine in this process might be obscure, the recognition accuracy of the classifier will be the judge of its quality of learning or its performance [1].

In some new classification systems, it is tried to investigate

the errors and propose a solution to compensate them [2-5]. There are many classification and clustering methods as well as the combinational approaches [6-8]. While the supervised learning tries to learn from the true labels or answers of the teacher, in semi-supervised the learner conversely uses teacher just to approve or not to approve the data in total. It means that in semi-supervised learning there is not really available teacher or supervisor. The procedure first starts with fully random manner, and when it reaches the state of final, it looks to the condition whether he won or lost. For example in the chess game, take it in consideration, that there may be none supervisor, but you gradually train to play better by trial-and-error process and looking at the end of the game to find you won or lost. K-Nearest Neighbor (KNN) classification is one of the most fundamental and simple classification methods. When there is little or no prior knowledge about the distribution of the data, the KNN method should be one of the first choices for classification. It is a powerful non-parametric classification system which bypasses the problem of probability densities completely [9]. The main idea is to classify an input query  $x$  into the most frequent tag in the set of its neighbor tags. The KNN rule classifies  $x$  by assigning it the label most frequently represented among the  $K$  nearest samples; this means that, a decision is made by examining the labels on the  $K$ -nearest neighbors and taking a vote. It is first introduced by Fix and Hodges in 1957 [10]. Later in 1967, KNN is looked at in theoretic perspective [11]. Once such consideration of KNN classification were established, a long line of investigation ensued including new rejection approaches [12], refinements with respect to Bayes error rate [13], distance weighted approaches [14, 15], soft computing [16] methods and fuzzy methods [17, 18].

ITQON et al. in [19] proposed a classifier, TFkNN, aiming at upgrading of distinction performance of KNN classifier and combining plural KNNs using testing characteristics. Their method not only upgrades distinction performance of the KNN but also brings an effect stabilizing variation of recognition ratio; and on recognition time, even when plural KNNs are performed in parallel, by devising its distance calculation it can be done not so as to extremely increase on comparison with that in single KNN. Some KNN advantages can be as follows: simplicity, robustness to noisy training data, and effectiveness in the adequate training data. It has some disadvantages such as: high computation cost in a test query, the large memory to implement, low accuracy rate in multidimensional data sets, parameter  $K$ , unclearness of distance type. Shall we use all attributes or certain attributes only [20]?

In this paper a new interesting algorithm is proposed which partially overcomes the low accuracy rate of KNN. Beforehand, it preprocesses the train set, computing the

validity of any train samples. Then the final classification is executed using weighted KNN which is employed the validity as the multiplication factor.

## II. MODIFIED K-NEAREST NEIGHBOR

The main idea of the presented method is assigning the class label of the queried instance into K validated data training points. In other hand, first, the validity of all data samples in the train set is computed. Then, a weighted KNN is performed on any test samples. Fig 1 shows the pseudo code of the MKNN algorithm.

```

Output_label := MKNN ( train_set , test_sample )
Begin
For i := 1 to train_size
Validity(i) := Compute Validity of i-th sample;
End for;
Output_label:=Weighted_KNN(Validity,test_sample);
Return Output_label ;
End.

```

Fig.1. Pseudo-code of the MKNN Algorithm

## III. VALIDITY OF THE TRAIN SAMPLES

In the MKNN algorithm, every training sample must be validated at the first step. The validity of each point is computed according to its neighbors. The validation process is performed for all train samples once. After assigning the validity of each train sample, it is used at the second step as impact or weight of the points in the ensembles of neighbors which the point is selected to attend.

To validate a sample point in the train set, the  $H$  nearest neighbors of the point is considered. Among the  $H$  nearest neighbors of a train sample  $x$ ,  $validity(x)$  counts the number of points with the same label to the label of  $x$ . Eq. 1 is the formula which is proposed to compute the validity of every points in train set.

$$Validity(x) = \frac{1}{H} \sum_{i=1}^H S(lbl(x), lbl(N_i(x))) \quad (1)$$

Where  $H$  is the number of considered neighbors and  $lbl(x)$  returns the true class label of the sample  $x$ . also,  $N_i(x)$  stands for the  $i$ th nearest neighbor of the point  $x$ . The function  $S$  takes into account the similarity between the point  $x$  and the  $i$ th nearest neighbor. Eq. 2 defines this function.

$$S(a, b) = \begin{cases} 1 & a = b \\ 0 & a \neq b \end{cases} \quad (2)$$

## IV. APPLYING WEIGHTED KNN

Weighted KNN is one of the variations of KNN method which uses the  $K$  nearest neighbors, regardless of their classes, but then uses weighted votes from each sample rather than a simple majority or plurality voting rule. Each of the  $K$  samples is given a weighted vote that is usually equal to some decreasing function of its distance from the unknown sample. For example, the vote might set be equal to  $1/(d_e+1)$ , where  $d_e$  is Euclidian distance. These weighted

votes are then summed for each class, and the class with the largest total vote is chosen. This distance weighted KNN technique is very similar to the window technique for estimating density functions. For example, using a weighted of  $1/(d_e+1)$  is equivalent to the window technique with a window function of  $1/(d_e+1)$  if  $K$  is chosen equal to the total number of training samples [21].

In the MKNN method, first the weight of each neighbor is computed using the  $1/(d_e+\alpha)$ , where  $\alpha$  is a smoothing regulator and here  $\alpha$  is selected to 0.5. Then, the validity of that training sample is multiplied on its raw weight which is based on the Euclidian distance. In the MKNN method, the weight of each neighbor sample is derived according to Eq. 3.

$$W(i) = Validity(i) \times \frac{1}{d_e + \alpha} \quad (3)$$

Where  $W(i)$  and  $Validity(i)$  stand for the weight and the validity of the  $i$ th nearest sample in the train set. This technique has the effect of giving greater importance to the reference samples that have greater validity and closeness to the test sample. So, the decision is less affected by reference samples which are not very stable in the feature space in comparison with other samples. In other hand, the multiplication of the validity measure on distance based measure can overcome the weakness of any distance based weights which have many problems in the case of outliers. So, the proposed MKNN algorithm is significantly stronger than the traditional KNN method which is based just on distance.

## V. EXPERIMENTAL RESULTS

This section discusses the experimental results and compares the MKNN method with original KNN algorithm.

### 1) Data sets

The proposed method is evaluated on nine standard data sets, namely Iris, Wine, Isodata, SAHeart, Balance-scale, Bupa and Monk's problems (including three problems). None of the databases had missing values, as well as they use continuous attributes. These standard data sets which are obtained from UCI repository [22] are described as follows. The iris database which is possibly one of the most frequently used benchmarks for evaluating classification and clustering algorithms is a well-defined problem with clear separating class boundaries. The data set contains 150 instances using three classes, where each class refers to a type of iris plant, namely Setosa, Versicolour and Virginica. This database uses four continuous attributes: sepal length, sepal width, petal length and petal width. The Wine data set is the result of a chemical analysis of wines grown in the same region in Italy but derived from three different cultivars. The analysis determined the quantities of 13 constituents found in each of the three types of wines. This data set has been used with many others for comparing various classifiers. In a classification context, this is a well-posed problem with well-behaved class structures. It has

three classes with 59, 71 and 48 instances. The more detail information about the wine data set is described in [23].

The Isodata set is the first test case in this study which is a two class data set and has 34 features as well as the 351 sample points. The SAHeart data set which is obtained from [www-stat.stanford.edu/ElemStatLearn](http://www-stat.stanford.edu/ElemStatLearn) is a retrospective sample of males in a heart-disease high-risk region of the Western Cape, South Africa. There are roughly two controls per case of CHD. Many of the CHD positive men have undergone blood pressure reduction treatment and other programs to reduce their risk factors after their CHD event. In some cases the measurements were made after these treatments. This data set has nine continuous features and two classes with the number of 463 instances. These data are taken from a larger dataset, described in [24].

The Balance-scale data set was generated to model psychological experimental results. Each example is classified as having the balance scale tip to the right, tip to the left, or be balanced which means three classes. The attributes are the left weight, the left distance, the right weight, and the right distance. It means this data set has four attributes. It has total 625 samples which include 49 balanced, 288 left, 288 right. The Bupa data set is a two class data set for classification. It contains 345 data sample as well as six attributes. In these six data sets, the instances are divided into training and test sets by randomly choosing 90% and 10% of instances per each of them, respectively. Also, all above mentioned data sets are become normalized with the mean of 0 and variance of 1,  $N(0,1)$  before applying the algorithms. The last experimented data set is Monk's problem which is the basis of a first international comparison of learning algorithms. There are three Monk's problems. The domains for all Monk's problems are the same. The second Monk's problem has added noise. For each problem, the domain has been partitioned into a train and test set. The number of Instances and attributes in all three problems are respectively, 432 and 6. These problems are two class problems. The train and test sets in all three Monk's problems are predetermined. The train sets in Monk 1, 2 and 3 are 124, 169 and 122, respectively.

## 2) Experiments

All experiments are evaluated over 500 independent runs and the average results of these examinations are reported. In all experiments, the number of considered neighbors (the value of parameter  $H$  in Eq. 1) is set to a fraction of the number of train data which is empirically set to 10% of the train size. Table 1 shows the results of the performance of classification using the presented method, MKNN, and traditional method, original version of KNN, comparatively. The experiments show that the MKNN method significantly outperforms the KNN method, with using different choices

of value  $K$ , over large variety of datasets. It is obvious that more information usually yields to more classification performance. Because of the MKNN classification is based on validated neighbors which have more information in comparison with simple class labels, it outperforms the KNN algorithm in performance. Fig 2 investigates the effect of parameter  $K$  on accuracy of algorithms KNN and MKNN comparatively in four different data sets: Iris, Balance-scale, Bupa and SAHeart. The value of  $K$  is the odd numbers in the range of [3-15]. Although usually the MKNN method initially overwhelms the KNN algorithm, the results of two algorithms gradually close to each other by growing the value of  $K$ . It can be because of the larger values of  $K$  result in invalidity of the validity of train samples. For some data sets the KNN even dominates the MKNN method with large  $K$  (see Fig 2b and 2c). In addition, since computing the validity measure is executed only once in training phase of the algorithm, computationally, the MKNN method can be applied with the nigh same burden in comparison with the weighted KNN algorithm.

## VI. CONCLUSION

In this paper, a new algorithm for improving the performance of KNN classifier is proposed which is called Modified K-Nearest Neighbor, MKNN. The proposed method which considerably improves the performance of KNN method employs a kind of preprocessing on train data. It adds a new value named "Validity" to train samples which cause to more information about the situation of training data samples in the feature space. The validity takes into accounts the value of stability and robustness of the any train samples regarding with its neighbors. Applying the weighted KNN which employs validity as the multiplication factor yields to more robust classification rather than simple KNN method, efficiently. The method evaluated on nine different benchmark tasks: Wine, Isodata, Iris, Bupa, Inosphere and three Monk's problems. The results confirm authors' claim about its robustness and accurateness unanimously. So this method is better in noisy datasets and also in the case of outliers. Since the outliers usually gain low value of validity, it considerably yields to robustness of the MKNN method facing with outliers. The experiments on Monk 2 approve this claim.

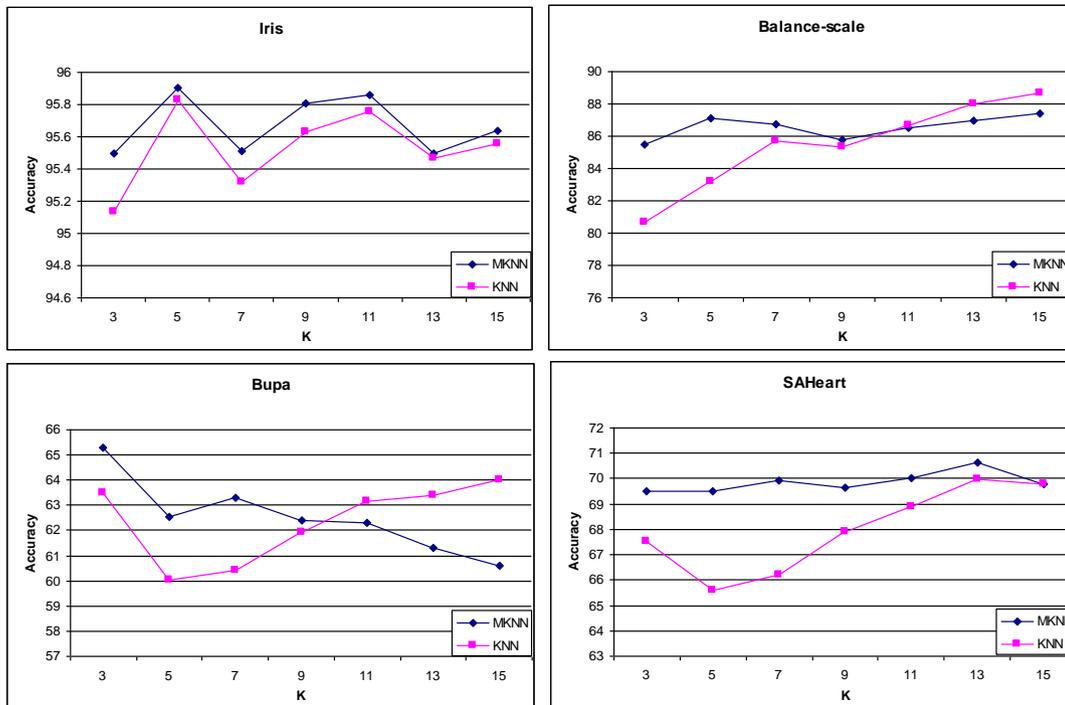


Fig.2. The effect of parameter K on accuracy of algorithms KNN and MKNN comparatively in four data sets (a) Iris (b) Balance scale (c) Bupa (d) SAHeart.

TABLE 1. Comparison of recognition rate between the MKNN and KNN algorithm (%).

	K=3		K=5		K=7	
	KNN	MKNN	KNN	MKNN	KNN	MKNN
<b>Monk 1</b>	84.49	<b>87.81</b>	84.26	<b>87.81</b>	79.86	<b>86.65</b>
<b>Monk 2</b>	69.21	<b>77.66</b>	69.91	<b>78.01</b>	65.74	<b>77.16</b>
<b>Monk 3</b>	89.12	<b>90.58</b>	89.35	<b>90.66</b>	88.66	<b>91.28</b>
<b>Isodata</b>	82.74	<b>83.52</b>	82.90	<b>83.32</b>	80.50	<b>83.14</b>
<b>Wine</b>	80.89	<b>83.95</b>	83.79	<b>85.76</b>	80.13	<b>82.54</b>
<b>Iris</b>	95.13	<b>95.50</b>	95.83	<b>95.90</b>	95.32	<b>95.51</b>
<b>Balance-sc</b>	80.69	<b>85.49</b>	83.22	<b>87.10</b>	85.74	<b>86.77</b>
<b>Bupa</b>	<b>63.51</b>	63.30	60.01	<b>62.52</b>	60.41	<b>63.29</b>
<b>SAHeart</b>	<b>67.51</b>	<b>69.51</b>	<b>65.59</b>	<b>69.49</b>	<b>66.21</b>	<b>69.95</b>

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