Liver Classification Using Modified Rotation Forest

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Abstract—Eesembling Classification techniques have been widely used in the medical field for accurate classification than an individual classifier. Modified Rotation Forest algorithm was proposed for accurate liver classification by analyzing the combination of selected classification algorithm and feature selection technique. Selected classification algorithms were considered from each category of classification algorithms. The category of classification algorithms are Tree based, Statistical based, Neural Networks based, Rule based and Lazy learners. Modified Rotation Forest algorithm for UCI liver data set has multi layer perception classification algorithm and Random Subset feature selection technique and for INDIA liver data set has nearest neighbour with generalized distance function and correlation based feature selection technique.

Keywords-Ensemble Classifier, Data Mining, Liver diagnosis, Classification, Feature selection

I. INTRODUCTION

Classification techniques are very popular in various automatic medical diagnoses tools. Problems with liver patients are not easily discovered in an early stage as it will be functioning normally even when it is partially damaged [1]. An early diagnosis of liver problems will increase patient's survival rate. Liver disease can be diagnosed by analyzing the levels of enzymes in the blood [].

Selected classification algorithms were combined with selected feature selection methods to identify the best combination of classification algorithm and feature selection to modify the Rotation Forest algorithm for accurate classification liver data.

In this paper, J48 and simple cart classification algorithms are from tree based algorithms, Naïve bayes and Bayes net classification algorithms are from statistical based algorithms, MLP and SMO classification algorithms are from multi layer perception based algorithms, IBK and KStar classification algorithms are from lazy learners and PART

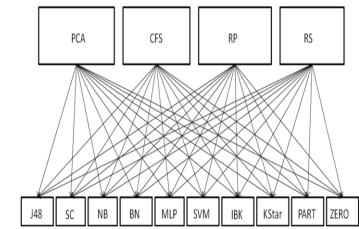


Figure 1. The combinations of classification algorithms and feature selection methods

and Zero classification algorithms are from rule based algorithms with all the combinations of Principal component analysis (PCA), Correlation based feature selection (CFS), Random projection and Random subset feature selection methods were considered for evaluating best combination of classification algorithm and feature selection method for modified rotation forest algorithm for UCI and INDIA liver data sets.

Two Liver patient data sets were used in this study, one is from Andhra Pradesh state of India and the second one is BUPA Liver Disorders datasets taken from University of California at Irvine (UCI) Machine Learning Repository [4]. In this experimentation, 10-fold cross-validation have been used. This paper concentrates on performance of combination of classification algorithms with feature selection methods.

II. RELATED WORK

Kun- Hong Liu and De-Shuang Huang [1] addressed the microarray dataset based cancer classification using rotation forest. Principal component analysis (PCA) was applied to feature transformation in the original rotation forest. In this paper Independent component analysis (ICA) was applied on breast cancer dataset and prostate dataset to validate the efficiency of rotation forest. The experimental results shows that ICA improves the performance of rotation forest compared with original transformation methods.

Juan J. Rodri'guez et al. [2] proposed a method for generating classifier ensembles based on feature extraction. The idea of the rotation approach is to encourage simultaneously individual accuracy and diversity within the ensemble. The experimental results with 33 data sets from UCI Machine Learning Repository showed that Rotation Forest outperformed all three methods by a large margin.

Akin Ozcift and Arif Gultenb [3] constructed rotation forest (RF) ensemble classifiers of 30 machine learning algorithms to evaluate their classification performances using Parkinson's, diabetes and heart diseases from literature. Experiments demonstrate that RF, as a newly proposed ensemble algorithm, proves itself to be efficient to increase classifier accuracies significantly.

Bendi Venkata Ramana et al. [5] compared popular Classification Algorithms for evaluating their classification performance in terms of Accuracy, Precision, Sensitivity and Specificity in classifying liver patients dataset. Accuracy, Precision, Sensitivity and Specificity are better for the AP Liver Dataset compared to UCLA liver datasets with all the selected algorithms. This can be attributed to more number of useful attributes like Total bilirubin, Direct bilirubin, Indirect bilirubin

Bendi Venkata Ramana et al. [7] proposed Bayesian Classification for diagnosis of liver diseases. The Bayesian Classification is combined with Bagging and Boosting for better accuracy. This accuracy can be further improved with huge amount of data.

Bendi Venkata Ramana et al. [8] proposed ANOVA and MANOVA for population comparision between ILPD data set and UCI data set. The results indicates that there exists more significant difference in the groups with all the possible attribute combinations except analysis on SGPT between non liver patients of UCI and INDIA data sets.

Bendi Venkata Ramana et al. [9] proposed Bayesian classification for accurate liver disease diagnosis and its accuracy further improved using

Wei Wanga et al. [11] introduced a multiple classifier ensemble approach based on performance evaluation, and apply it to segmenting tissue regions, especially the biomarker aceto white tissue in digitized uterine cervix images. The multiple classifier system uses a multi-observer segmentation evaluation tool (MOSES) to train and combine SVM classifiers. Experimental results show that the proposed classifier ensemble performs better than a single SVM classifier.

Bing Chen and Hua-Xiang Zhang [12] proposed an approach of multiple classifiers ensemble based on feature selection (FSCE) In order to improve the classification performance of classifiers. FSCE is tested on the UCI benchmark data sets, and compared classification efficiency with member classifiers trained based on the algorithm of Adaboost.

Manju Bhardwaj et al. [13] proposed and empirically evaluate a novel method for generating members of ensemble based on 'learning from- mistakes' paradigm. SVM is used as the base learner, and a series of dependent classifiers is obtained using model based instance selection method. Simple majority voting has been used to combine learners. It is found that the ensemble created shows better accuracy as compared to the ensembles created using AdaBoost, MAdaBoost, Bagging and Arc-x4SVM.

Lishuang LI [14] presented a gene mention tagging system which utilizes the two-layer stacking classifiers ensemble method to combine multiple powerful machine learning frameworks. The two-layer stacking algorithm is a more effective classifiers ensemble method than voting and simple set operation methods.

M. F. Amasyali [15] compared 12 single classifiers and 11 classifier ensembles over 36 datasets according to Classification accuracy and execution time. According to classification accuracy the best 6 algorithms are ordered as Rotation Forest, Random Committees, Random Forest, Logit Boost, Decorate and Bagging. The results show that Rotation Forest has the highest accuracy. However, when accuracy and execution time are considered together, Random Forest and Random Committees can be the best choices.

Aleksandar Lazarevic and Zoran Obradovic[16] proposed several methods for pruning neural network ensembles. The clustering based approach applies k-means clustering to entire set of classifiers in order to identify the groups of similar classifiers and then eliminates redundant classifiers inside each cluster. The novel proposed methods applied to several data sets have shown that by selecting an optimal subset of neural network classifiers, it is possible to obtain significantly smaller ensemble of classifiers while achieving the same or even slightly better generalizability as when using the entire ensemble.

Jin Zhou [17] proposed a new ensemble method for cancer data classification. The gene expression data is preprocessed for normalization and then Kernel Independent Component Analysis (KICA) is applied to extract features. Then an intelligent approach is brought forward, which uses Support Vector Machine (SVM) as the base classifier and applied with Binary Particle Swarm Optimization (BPSO) for constructing ensemble classifiers. The leukemia and colon datasets are used for conducting all the experiments. Results show that the proposed method produces a good recognition rate comparing with some other advanced artificial techniques.

Nigar Sen Koktas [18] discussed ensemble methods for gait classification as a part of preliminary studies of designing a semi-automated diagnosis system. For this purpose gait data is collected from 110 sick subjects (having knee Osteoarthritis (OA)) and 91 age-matched normal subjects. A set of Multilayer Perceptions (MLPs) is trained by using joint angle and time-distance parameters of gait as features. The result suggests that ensemble of MLPs could be applied in the automated diagnosis of gait disorders in a clinical context.

Yun Zhai [19] proposed an ensemble classifier model SCECM for imbalanced datasets in a heterogeneous classifier fusion environment, which is based on a differentiated re-sampling technology DSRA and an improved Adaboost algorithmComparative experimental results show that the proposed approach improves performance for the minority class while preserving the ability to recognize examples from the majority classes.

Hualong Yu [20] proposed a novel ensemble classification method named as simple rule-based ensemble classifiers (SREC) and used it for cancer micro array data classification. Experimental results indicate the proposed method is effective and feasible because it has produced less classification errors than many other classifiers. Meanwhile, it has some other advantages such as low time-complexity and storage space etc.

Sung-Bae Cho [21] presented a method to search optimal ensemble of diverse feature-classifier pairs using speciated FA. This paper proposes sophisticated ensemble of such features and classifiers to obtain high classification performance. Speciated genetic algorithm has been exploited to get the diverse ensembles of features and classifiers in a reasonable time. Experimental results with two well-known datasets indicate that the proposed method finds many good ensembles that are superior to other individual classifiers.

III. CLASSIFICATION ALGORITHMS

The classification techniques are categorized into Eager learners, Lazy learners (learning from neighbors) and other classification techniques. Eager learners will take a set of training tuples, and construct a classification model (learned model) and eager to classify previously unseen tuples. The classification methods belongs to eager learners are decision tree induction, Bayesian classification, rule-based classification, classification by back propagation, support vector machines, and classification based on association rule mining. The Eager learners are tree based classifiers, statistical based classifiers, Rule based classifiers and neural network based classifiers and K-nearest-neighbor classifier is a Lazy learner.

A. Tree Based Algorithms

Decision tree induction is the learning of decision trees from class-labeled training tuples. A decision tree is a flowchart-like tree structure, where each internal node (non leaf node) denotes a test on an attribute, each branch represents an outcome of the test, and each leaf node (or terminal node) holds a class label. The topmost node in a tree is the root node.

Check for base cases For each attribute a Find the feature that best divides the training data such as information gain from splitting on a Let a best be the attribute with the highest normalized information gain Create a decision node node that splits on a_best Recurse on the sub-lists obtained by splitting on a_best and add those nodes as children of node *Figure 2.* Algorithm for building a decision tree

B. Statistical based algorithms

Statistical classifiers based on bayes theorem. Statistical classifiers shows high accuracy and speed when applied to large data bases. It works on one assumption called class conditional independence that is the effect of an attribute value on a given class is independent of the values of the other attributes.

```
Procedure Bayesian_Classifier (X = <X1,,,Xn > ):

Begin

For all classes Ci \in C = C1, , , Cm

{

Compute P(Ci):

For all features Xj \in X

{

Compute P (Xj / Ci):

};

p(\chi_j / C_i)'s(=\prod_{j=1}^n p(\chi_j / C_i));

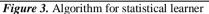
Multiply all

f_i(d) = p(C_i)^* \prod_{j=1}^n p(\chi_j / C_i));

Calculate

};

Assign d to the class (es) of max (f1 (d),..., fm(d)) End
```



C. Neural network based algorithms

A multi-layer neural network consists of large number of units (neurons) joined together in a pattern of connections. Units in a net are usually segregated into three classes: input units, which receive information to be processed; output units, where the results of the processing are found; and units in between known as hidden units. Feed-forward ANNs allow signals to travel one way only, from input to output.

1. Present a training sample to the neural network.

Compare the network's output to the desired output from that sample. Calculate the error in each output neuron.
 For each neuron, calculate what the output should have been, and a scaling factor, how much lower or higher the output must be adjusted to match the desired output. This is the local error.
 Adjust the weights of each neuron to lower the local error.
 Assign "blame" for the local error to neurons at the previous level, giving greater responsibility to neurons connected by stronger weights.
 Repeat the steps above on the neurons at the previous level, using each one's "blame" as its error.
 With more details, the general rule for updating weights is:

$$\Delta w_{ji} = \eta \delta_j o_i;$$

where:

• η is a positive number (called learning rate), which determines the step size in the gradient descent search. A large value enables back propagation to move faster to the target weight configuration but it also increases the chance of its never reaching this target.

• Oi is the output computed by neuron i

$$\partial_i = O_i(1 - O_i)(T_i - O_i);$$

for the output neurons, where Ti the wanted output for the neuron i and

$$\delta_j = o_j (1 - o_j) \sum_k \delta_k w_{kj};$$

for the internal (hidden) neurons

Figure 4. Algorithm for neural network

D. Lazy learners

Lazy-learning algorithms require less computation time during the training phase than eager-learning algorithms (such as decision trees, neural and Bayes nets) but more computation time during the classification process. One of the most straightforward lazy learning algorithms is the nearest neighbour algorithm.

Procedure Lazy Learner (Testing Instances)						
for	each	testing	instance			
	tes of the training set according to class label of the k nearest instances	a distance metric				

Figure 5. Algorithm for lazy learners

E. Rule based algorithms

Decision trees can be translated into a set of rules by creating a separate rule for each path from the root to a leaf in the tree (Quinlan, 1993). However, rules can also be directly induced from training data using a variety of rulebased algorithms. Furnkranz (1999) provided an excellent overview of existing work in rule-based methods.

On presentation of training examples							
Training exam	Training examples:						
Initialize rule	Initialize rule set to a default (usually empty, or a rule assigning all objects to the most common class).						
Initialize examples to either all available examples or all examples not correctly handled by rule set.							
Repeat	Repeat						
(a) Find best, the best rule with respect to examples.							
(b)	If	such	а	rule	can	be	found
i. Add best to rule set.							
ii. Set examples to all examples not handled correctly by rule set.							
until no rule best can be found (for instance, because no examples remain).							
Figure 6 Algorithm for rule learners							

Figure 6. Algorithm for rule learners

F. Support Vector Machines

Support Vector Machines (SVMs) are the newest supervised machine learning technique (Vapnik, 1995). SVMs revolve around the notion of a "margin "either side of a hyper plane that separates two data classes. Maximizing the margin and thereby creating the largest possible distance between the separating hyper plane and the instances on either side of it has been proven to reduce an upper bound on the expected generalization error.

Introduce positive Lagrange multipliers, one for each of the inequality constraints This gives Lagrangian: $L_p = \frac{1}{2} || w ||^2 - \sum_{i=1}^{N} \alpha_i y_i (x_i \cdot w - b) + \sum_{i=1}^{N} \alpha_i$ Minimize LP with respect to w, b.

This is a convex quadratic programming problem. In the solution, those points for which α i >0 are called "support vectors"

Figure 7. Algorithm for SVMs

IV. RESULTS AND DISCUSSION

In this study we used liver data sets from UCI and INDIA. UCI liver data set has 345 samples with 6 features and INDIA liver data set has 583 samples with 10 features. For the purpose of experimentation, Weka© Data Mining open source machine learning software [10]. It is used on i7 processor with 4 GB RAM. With Each combination of classification algorithm and feature selection method, we have observed Accuracy which can be defined as follows: Accuracy: The accuracy of a classifier is the percentage of the test set tuples that are correctly classified by the classifier.

$$Accuracy = \frac{IP + IN}{TP + FP + FN + TN}$$

TD TN

Where TP means True Positives, TN means True Negatives, FP means False Positives and FN means False Negatives

Table 1: Liver Patient data set and attributes					
Attribute	Туре				
Gender	Categorical				
Age	Real number				
Total_bilirubin	Real number				
Direct_ bilirubin	Real number				
Total_protiens	Real number				
Albumin	Real number				
A/G ratio	Real number				
SGPT	Integer				
SGOT	Integer				
ALP	Integer				

Table 2: UCLA Liver data set and attributes available

Attribute	Туре
Mcv	Integer
Alkphos	Integer
SGPT	Integer
SGOT	Integer
Gammagt	Real number

The accuracy of all combinations of ten classification algorithms which were selected from various categories of classification algorithms and feature selection techniques were analyzed for UCI liver data set and INDIA liver data set.

Feature Selection		PCA	CFS	Random	Random
				Projection	Subset
	J48	73.3333	63.7681	67.2464	68.1159
	Simple Cart	73.3333	66.087	68.9855	69.5652
smi	Naive Bayes	54.7826	64.058	56.8116	56.8116
Classification Algorithms	Bayes Net	71.3043	57.971	56.2319	56.5217
	MLP	71.3043	63.7681	68.4058	74.7826
	SMO	58.5507	57.971	57.971	57.971
tiio	IBK	63.1884	64.058	66.9565	61.1594
Issifica	KStar	67.5362	68.4058	66.6667	67.5362
	PART	72.4638	63.4783	68.1159	69.2754
Cla	Zero	57.971	57.971	57.971	57.971

Table 3: UCI Liver Data Set

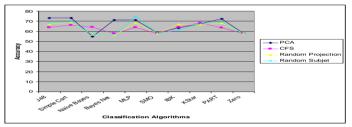


Figure 8. Accuracy for different combination of classifier and feature selection for UCI liver data set

Table 4: INDIA Liver Data Set						
Feature Selection		PCA	CFS	Random	Random	
				Projection	Subset	
	J48	69.8113	71.012	70.669	69.2967	
	Simple Cart	69.9828	71.3551	71.012	70.4974	
smi	Naive Bayes	54.2024	56.2607	55.2316	55.06	
nith	Bayes Net	64.1509	68.9537	65.0086	68.4391	
lgc	MLP	70.1544	71.3551	71.5266	71.5266	
ЧЧ	SMO	71.3551	71.3551	71.3551	71.3551	
tio	IBK	68.6106	69.2967	71.3551	69.6398	
Classification Algorithms	KStar	71.8696	73.0703	70.669	71.3551	
	PART	70.3259	71.012	70.3259	68.4391	
Cla	Zero	71.3551	71.3551	71.3551	71.3551	

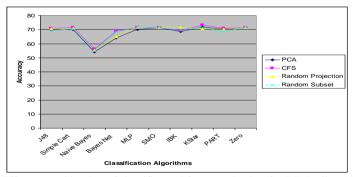


Figure 9. Accuracy for different combination of classifier and feature selection for INDIA liver data set

Training Phase

Given X: the objects in the training data set (an N X n matrix) Y: the labels of the training set (an N X 1 matrix) L: the number of classifiers in the ensemble K: the number of subsets

{ $\mathcal{O}_{1,\ldots,\mathcal{O}_{2}}$: the set of class lables

For i=1...L Prepare the rotation matrix R_i^a Split F(the feature set) into k subsets: F_{ij} (for j=1...k) For j=1...k Let X_{ij} be the data set X for the features in F_{ij} Eliminate from X_{ij} a random subset of classes Select a bootstrap sample from X_{ij} of size 75% of the number of objects in X_{ij} . Denote the new set by X'_{ij} Apply Random subset on X'_{ij} to obtain the coefficients in matrix C_{ij} Arrange the C_{ij} for j=1...k in a rotation matrix R_i^a as in equation (1) Construct R_i^a by rearranging the columns of R_i so as to match the order of features in F Build classifier D_i using (X R_i^a , Y) as the training set

Classification Phase

For a given X, let $d_{i,j}$ (x R_i^a) be the probability assigned by the classifier D_i to the hypothesis that x comes from class \mathcal{O} j. Calculate the confidence for each class \mathcal{O} j, by the average combination method:

$$\mu_j(x) = \frac{1}{L} \sum_{i=1}^{L} d_{i,j}(xR^a_i), j = 1, \dots, c.$$

Assign x to the class with the largest confidence.

Figure 10. Modified rotation forest algorithm with random subset feature selection and multi layer perception classification algorithm for UCI liver data set

Training Phase Given

X: the objects in the training data set (an N X n matrix) Y: the labels of the training set (an N X 1 matrix) L: the number of classifiers in the ensemble K: the number of subsets { \emptyset 1,..., \emptyset 2}: the set of class lables For i=1...L Prepare the rotation matrix R_i^a Split F(the feature set) into k subsets: F_{ii} (for j=1...k) For j=1...k Let X_{ij} be the data set X for the features in F_{ii} Eliminate from X_{ii} a random subset of classes Select a bootstrap sample from X_{ii} of size 75% of the number of objects in X_{ii} . Denote the new set by X'_{ii} Apply CFS on $\hat{X'}_{ij}$ to obtain the coefficients in matrix C_{ij} Arrange the C_{ii} for j=1...k in a rotation matrix R_i^a as in equation (1)Construct R_i^a by rearranging the columns of R_i so as to match the order of features in F Build classifier D_i using (X R_i^a , Y) as the training set

Classification Phase

For a given X, let $d_{i,j}$ (x R_i^a) be the probability assigned by the classifier D_i to the hypothesis that x comes from class \mathcal{O} j. Calculate the confidence for each class \mathcal{O} j, by the average combination method:

$$\mu_j(x) = \frac{1}{L} \sum_{i=1}^{L} d_{i,j}(xR^a_i), j = 1, \dots, c.$$

Assign x to the class with the largest confidence.

Figure 11. Modified rotation forest algorithm with CFS feature selection and nearest neighbor classification algorithm for INDIA liver data set

The results shows that multi layer perception classification algorithm with random subset gives highest accuracy that is 74.7826 for the UCI liver data set and nearest neighbor with CFS gives highest accuracy that is 73.0703 for the INDIA liver data set.

V. CONCLUSIONS

In this study, ten popular Classification Algorithms were considered with all the combinations of four feature selection methods for evaluating their classification performance in terms of accuracy in classifying liver patient data sets. Modified rotation forest algorithm was proposed with multi layer perception classification algorithm and random subset feature selection method for UCI liver data set. Modified rotation forest algorithm was proposed with nearest neighbor classification algorithm and correlation based feature selection method for INDIA liver data set.

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