

EVOLUTIONARY COMPUTATION APPROACH TO ECG SIGNAL CLASSIFICATION

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Abstract: In this paper, we propose a novel classification system for ECG signals based on particle swarm optimization (PSO). The main objective of this system is to optimize the performance of the support vector machine (SVM) classifier in terms of accuracy by automatically: i) searching for the best subset of features where to carry out the classification task; and ii) solving the SVM model selection issue. Experiments conducted on the basis of ECG data from the MIT-BIH arrhythmia database to classify five kinds of abnormal waveforms and normal beats confirm the effectiveness of the proposed PSO-SVM classification system.

1 INTRODUCTION

The recent literature reports different and interesting methodologies for the automatic classification of electrocardiogram (ECG) signals (e.g., de Chazal et Reilly, 2006, and Inan et Giovangrandi, 2006). However, in the design of an ECG classification system, there are still some open issues, which if suitably addressed may lead to the development of more robust and efficient classifiers. One of these issues is related to the choice of the classification approach to be adopted. In particular, we think that, despite its great potential, the SVM approach has not received the attention it deserves in the ECG classification literature compared to other research fields. Indeed, the SVM classifier exhibits a promising generalization capability thanks to the maximal margin principle (MMP) it is based on (Vapnik, 1998). Another important property is its lower sensitivity to the curse of dimensionality compared to traditional classification approaches. This is explained by the fact that the MMP makes unnecessary to estimate explicitly the statistical distributions of classes in the hyperdimensional feature space in order to carry out the classification task. Thanks to these interesting properties, the SVM classifier proved successful in numerous and

different application fields, such as 3D object recognition (Pontil et Verri, 1998), biomedical imaging (El-Naqa et al., 2002), remote sensing (Melgani et Bruzzone, 2004 and Bazi et Melgani, 2006). Turning back to ECG classification, other issues which can be identified are: 1) feature selection is not performed in a completely automatic way; and 2) the selection of the best free parameters of the adopted classifier is generally performed empirically (model selection issue).

In this paper, we present in a first step a thorough experimental exploration of the SVM capabilities for ECG classification. In a second step, in order to address the aforementioned issues, we propose to optimize further the performances of the SVM approach in terms of classification accuracy by 1) automatically detecting the best discriminating features from the whole considered feature space and 2) solving the model selection issue. The detection process is implemented through a particle swarm optimization (PSO) framework that exploits a criterion intrinsically related to the SVM classifier properties, namely the number of support vectors (#SV).

2 PROPOSED APPROACH

2.1 Support Vector Machine (SVM)

Let us first for simplicity consider a supervised binary classification problem. Let us assume that the training set consists of N vectors $\mathbf{x}_i \in \mathcal{R}^d$ ($i = 1, 2, \dots, N$) from the d -dimensional feature space \mathbf{X} . To each vector \mathbf{x}_i , we associate a target $y_i \in \{-1, +1\}$. The linear SVM classification approach consists of looking for a separation between the two classes in \mathbf{X} by means of an optimal hyperplane that maximizes the separating margin (Vapnik, 1998). In the nonlinear case, which is the most commonly used as data are often linearly nonseparable, they are first mapped with a kernel method in a higher dimensional feature space, i.e., $\Phi(\mathbf{X}) \in \mathcal{R}^{d'}$ ($d' > d$). The membership decision rule is based on the function $\text{sign}[f(\mathbf{x})]$, where $f(\mathbf{x})$ represents the discriminant function associated with the hyperplane in the transformed space and is defined as:

$$f(\mathbf{x}) = \mathbf{w}^* \cdot \Phi(\mathbf{x}) + b^* \quad (1)$$

The optimal hyperplane defined by the weight vector $\mathbf{w}^* \in \mathcal{R}^{d'}$ and the bias $b^* \in \mathcal{R}$ is the one that minimizes a cost function that expresses a combination of two criteria: margin maximization and error minimization. It is expressed as:

$$\Psi(\mathbf{w}, \xi) = \frac{1}{2} \|\mathbf{w}\|^2 + C \sum_{i=1}^N \xi_i \quad (2)$$

This cost function minimization is subject to the following constraints:

$$y_i (\mathbf{w} \cdot \Phi(\mathbf{x}_i) + b) \geq 1 - \xi_i, \quad i = 1, \dots, N \quad (3)$$

and

$$\xi_i \geq 0, \quad i = 1, 2, \dots, N \quad (4)$$

where the ξ_i 's are slack variables introduced to account for non-separable data. The constant C represents a regularization parameter that allows to control the shape of the discriminant function. The above optimization problem can be reformulated through a Lagrange functional, for which the Lagrange multipliers can be found by means of a dual optimization leading to a Quadratic Programming (QP) solution, i.e.,

$$\max_{\alpha} \sum_{i=1}^N \alpha_i - \frac{1}{2} \sum_{i,j=1}^N \alpha_i \alpha_j y_i y_j K(\mathbf{x}_i, \mathbf{x}_j) \quad (5)$$

under the constraints:

$$\alpha_i \geq 0 \quad \text{for } i = 1, 2, \dots, N \quad (6)$$

and

$$\sum_{i=1}^N \alpha_i y_i = 0 \quad (7)$$

where $\alpha = [\alpha_1, \alpha_2, \dots, \alpha_N]$ is the vector of Lagrange multipliers and $K(\cdot, \cdot)$ is a kernel function. The final result is a discriminant function conveniently expressed as a function of the data in the original (lower) dimensional feature space \mathbf{X} :

$$f(\mathbf{x}) = \sum_{i \in S} \alpha_i^* y_i K(\mathbf{x}_i, \mathbf{x}) + b^* \quad (8)$$

The set S is a subset of the indices $\{1, 2, \dots, N\}$ corresponding to the non-zero Lagrange multipliers α_i 's, which define the so-called support vectors.

As described above, SVMs are intrinsically binary classifiers. But the classification of ECG signals often involves the simultaneous discrimination of numerous information classes. In order to face this issue, different multiclass classification strategies can be adopted (Melgani et Bruzzone, 2004). In this paper, we shall consider the commonly used one-against-all strategy.

2.2 PSO Principles

Particle swarm optimization (PSO) is a stochastic optimization technique introduced recently by Kennedy and Eberhart, which is inspired by social behavior of bird flocking and fish schooling (Kennedy et Eberhart, 2001). Similarly to other evolutionary computation algorithms such as genetic algorithms (GAs) (Bazi et Melgani, 2006), PSO is a population-based search method, which exploits the concept of social sharing of information. This means that each individual (called *particle*) of a given population (called *swarm*) can profit from the previous experiences of all other individuals from the same population. During the iterative search process in the d -dimensional solution space, each particle (i.e., candidate solution) will adjust its flying velocity and position according to its own flying experience as well as the experiences of the other companion particles of the swarm.

Let us consider a swarm of size S . Each particle P_i ($i = 1, 2, \dots, S$) from the swarm is

characterized by: 1) its current position $\mathbf{p}_i(t) \in \mathfrak{R}^d$, which refers to a candidate solution of the optimization problem at iteration t ; 2) its velocity $\mathbf{v}_i(t) \in \mathfrak{R}^d$; and 3) the best position $\mathbf{p}_{bi}(t) \in \mathfrak{R}^d$ identified during its past trajectory. Let $\mathbf{p}_g(t) \in \mathfrak{R}^d$ be the best global position found over all trajectories traveled by the particles of the swarm. The position optimality is measured by means of one or more fitness functions defined in relation to the considered optimization problem. During the search process, the particles move according to the following equations:

$$\begin{cases} \mathbf{v}_i(t+1) = w\mathbf{v}_i(t) + c_1 \cdot r_1(t)(\mathbf{p}_{bi}(t) - \mathbf{p}_i(t)) \\ \quad + c_2 \cdot r_2(t)(\mathbf{p}_g(t) - \mathbf{p}_i(t)) \\ \mathbf{p}_i(t+1) = \mathbf{p}_i(t) + \mathbf{v}_i(t) \end{cases} \quad (9)$$

where $r_1(\cdot)$ and $r_2(\cdot)$ are random variables drawn from a uniform distribution in the range $[0, 1]$ so that to provide a stochastic weighting of the different components participating in the particle velocity definition. c_1 and c_2 are two acceleration constants regulating the relative velocities with respect to the best global and local positions, respectively. The inertia weight w is used as a tradeoff between global and local exploration capabilities of the swarm.

2.3 PSO Setup

The position $\mathbf{p}_i \in \mathfrak{R}^{d+2}$ of each particle P_i from the swarm is viewed as a vector encoding: 1) a candidate subset F of features among the d available input features, and 2) the value of the two SVM classifier parameters, which are the regularization and the kernel parameters C and γ , respectively. Since the first part of the position vector implements a feature detection task, each component (coordinate) of this part will assume either a "0" (feature discarded) or a "1" (feature selected) value. The conversion from real to binary values will be done by a simple thresholding operation at the 0.5 value.

Let $f(i)$ be the fitness function value associated with the i th particle P_i . The choice of the fitness function is important since, on its basis, the PSO evaluates the goodness of each candidate solution \mathbf{p}_i for designing our SVM classification system. A possible choice is to adopt the class of criteria that estimates the leave-one-out error bound, which exhibits the interesting property of representing an unbiased estimation of the generalization performance of classifiers. In particular, for SVM classifiers, different measures of this error bound have been derived, such as the radius-margin bound and the simple support vector (SV) count (Vapnik,

1998). In this paper, we will explore the simple SV count as fitness criterion in the PSO optimization framework because of its simplicity and effectiveness as shown in the context of the classification of hyperspectral remote sensing images (Bazi et Melgani, 2006).

2.4 SVM Classification with PSO

- **Initialization**

Step 1: Generate randomly an initial swarm of size S .

Step 2: Set to zero the velocity vectors \mathbf{v}_i ($i=1, 2, \dots, S$) associated with the S particles.

Step 3: For each position $\mathbf{p}_i \in \mathfrak{R}^{d+2}$ of the particle P_i ($i=1, 2, \dots, S$) from the swarm, train an SVM classifier and compute the corresponding fitness function $f(i)$ (i.e., the #SV measure).

Step 4: Set the best position of each particle with its initial position, i.e.,

$$\mathbf{p}_{bi} = \mathbf{p}_i, (i=1, 2, \dots, S) \quad (11)$$

- **Search process**

Step 5: Detect the best global position \mathbf{p}_g in the swarm exhibiting the minimal value of the considered fitness function over all explored trajectories.

Step 6: Update the speed of each particle using Equation (9).

Step 7: Update the position of each particle using Equation (10). If a particle goes beyond the predefined boundaries of the search space, truncate the updating by setting the position of the particle at the space boundary and reverse its search direction (i.e., multiply its speed vector by -1). This will permit to forbid the particles from further attempting to go outside the allowed search space.

Step 8: For each candidate particle \mathbf{p}_i ($i=1, 2, \dots, S$), train an SVM classifier and compute the corresponding fitness function.

Step 9: Update the best position \mathbf{p}_{bi} of each particle if its new current position \mathbf{p}_i ($i=1, 2, \dots, S$) has a smaller fitness function.

- **Convergence**

Step 10: If the maximal number of iterations is not yet reached, return to *Step 5*.

- **Classification**

Step 11: Select the best global position \mathbf{p}_g^* in the swarm and train an SVM classifier fed with the subset of detected features mapped by \mathbf{p}_g^* and modeled with the values of the two parameters C and γ encoded in the same position.

Step 12: Classify the ECG signals with the trained SVM classifier.

3 EXPERIMENTAL RESULTS

Our experiments were conducted on the basis of ECG data from the MIT-BIH arrhythmia database (Mark et Moody, 1997). In particular, the considered beats make reference to the following classes: normal sinus rhythm (N), atrial premature beat (A), ventricular premature beat (V), right bundle branch block (RB), left bundle branch block (LB), and paced beat (/). Similarly to (Inan et al., 2006), the beats were selected from the recordings of 18 patients, which correspond to the following files: 100, 102, 104, 105, 106, 107, 118, 119, 200, 201, 203, 205, 208, 212, 213, 214, 215, and 217. For feeding the classification process, we adopted in this study the two following kinds of features: i) ECG morphology features; and ii) three ECG temporal features that are the QRS complex duration, the RR interval (i.e., time span between two consecutive R points representing the distance between the QRS peaks of the present and previous beats), and the RR interval averaged over the ten last beats (de Chazal et Reilly, 2006). The total number of morphology and temporal features is equal to 303 for each beat. In order to train the classification process and to assess its accuracy, we selected randomly from the considered recordings 500 beats for the training set, whereas 42185 beats were used as test set (thus, the training set represents just 1.18% of the test set). The detailed numbers of training and test beats are reported for each class in Table 1. Classification performance was evaluated in terms of three accuracy measures, which are: 1) the overall accuracy (OA); 2) the accuracy of each class; and 3) the average accuracy (AA).

Due to the good performances generally achieved by the nonlinear SVM classifier based on the Gaussian kernel [6], we adopted this kernel in all experiments. The parameters C and γ were varied in the ranges $[10^{-3}, 200]$ and $[10^{-3}, 2]$, respectively. The k value and the number of hidden nodes (h) of the kNN and the RBF classifiers were tuned in the intervals $[1, 15]$ and $[10, 60]$, respectively. Concerning the PSO algorithm, we considered the following standard parameters: swarm size $S=40$, inertia weight $w=0.4$, acceleration constants c_1 and c_2 equal to the unity, and maximum number of iterations fixed to 40.

3.1 Experiment 1: Classification in the Original Feature Space

In this experiment, we applied the SVM classifier directly on the whole original hyperdimensional

feature space which is composed of 303 features. During the training phase, the SVM parameters (i.e., C and γ) were selected according to a m -fold cross-validation (CV) procedure. In all experiments reported in this paper, we adopted a 5-fold CV. The same procedure was adopted to find the best parameters for the kNN and the RBF classifiers. The best values obtained for the three investigated classifiers are $C=25$, $\gamma=0.5$, $k=3$ and $h=20$. As reported in Table 2, the OA and AA accuracies achieved by the SVM classifier on the test set are equal to 87.95% and 87.60%, respectively. These results are much better than those achieved by the RBF and the kNN classifiers. Indeed, the OA and AA accuracies are equal to 82.78% and 82.34% for the RBF classifier, and 78.21% and 79.34% for the kNN classifier, respectively. This experiment appears to confirm what observed in other application fields, i.e., the superiority of SVM with respect to traditional classifiers when dealing with feature spaces of very high dimensionality.

3.2 Experiment 2: Classification based on Feature Reduction

In this experiment, we trained the SVM classifier in feature subspaces of various dimensionalities. The desired number of features was varied from 10 to 50 with a step of 10, namely from small to high dimensional feature subspaces. Feature reduction was achieved by means of the traditional Principal Component Analysis (PCA) algorithm. Figure 1-a depicts the results obtained in terms of OA by the three considered classifiers combined with the PCA algorithm, namely the PCA-SVM, the PCA-RBF and the PCA-kNN classifiers. In particular, it can be seen that, for all feature subspace dimensionalities except the lowest one (i.e., 10 features), the PCA-SVM classifier maintains a clear superiority over the two other classifiers. Its best accuracy was found by using a feature subspace composed of the first 40 components. The corresponding OA and AA accuracies are equal to 88.98% and 88%, respectively. Comparing these results with those obtained by the SVM classifier in the original feature space (i.e., without feature reduction), a slight increase of 1.03% and 0.4% in terms of OA and AA, respectively, was achieved (see Table 2).

3.3 Experiment 3: Classification with PSO-SVM

In this experiment, we applied the PSO-SVM classifier on the available training beats. At

convergence of the optimization process, we assessed the PSO-SVM classifier accuracy on the test samples. The achieved overall and average accuracies are equal to 91.44% and 91.19% corresponding to substantial accuracy gains with respect to what yielded either by the SVM classifier applied on all available features (+3.49% and +3.59%, respectively) or by the PCA-SVM classifier (+2.46% and +3.19%, respectively) (see Table 2 and Figure 1). Its worst class accuracy was obtained for atrial premature beats (A) (88.16%) while that of the SVM and the PCA-SVM classifiers corresponded to paced beats (/) (73.43%) and ventricular premature beats (V) (78.06%), respectively. This shows the capability of the PSO-SVM classifier to reduce significantly the gap between the worst and best class accuracies (8.25% against 15.43% and 20.21% for the PCA-SVM and the SVM classifiers, respectively) while keeping the overall accuracy to a high level.

4 CONCLUSIONS

The main novelty of this paper is to be found in the proposed PSO-based approach that aims at optimizing the performances of SVM classifiers in terms of classification accuracy by detecting the best subset of available features and by solving the tricky model selection issue. Its completely automatic nature renders it particularly useful and attractive. The results confirm that the PSO-SVM classification system boosts up significantly the generalization capability achievable with the SVM classifier. Finally, it is noteworthy that the general nature of the proposed PSO-SVM system makes it applicable not just to morphology and temporal features but to other kinds of features such as those based on wavelets and high-order statistics. Finally, other optimization criteria could be considered as well individually or jointly depending on the application requirements.

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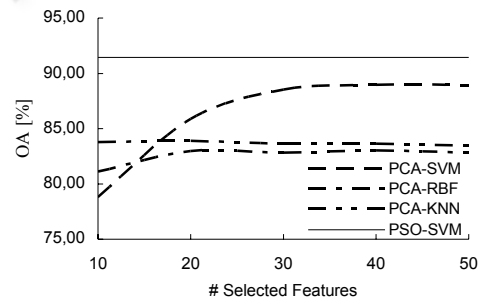


Figure 1: Overall percentage accuracy (OA) versus number of selected features (first principal components) achieved on the test beats by the different classifiers.

Table 1: Numbers of training and test beats used in the experiments.

Class	N	A	V	RB	/	LB	Total
Training beats	150	100	100	50	50	50	500
Test beats	24966	119	4239	3939	6971	1951	42185

Table 2: Overall (OA), average (AA), and class percentage accuracies achieved on the test beats by the different investigated classifiers.

Method	OA	AA	N	A	V	RB	/	LB
SVM	87.95	87.60	90.05	83.19	92.12	93.15	73.43	93.64
RBF	82.78	82.34	85.14	78.99	90.39	86.74	66.53	86.26
kNN	78.21	79.34	76.50	66.38	71.99	93.27	75.92	92.00
PCA-SVM	88.98	88.00	89.36	83.19	78.06	93.50	90.60	93.28
PCA-RBF	83.04	82.11	85.86	80.67	87.85	83.87	68.85	85.54
PCA-kNN	83.91	82.02	85.62	69.74	79.05	93.04	73.89	90.77
PSO-SVM	91.44	91.19	91.12	88.16	93.70	93.70	92.01	96.41

