Chapter 1

Controlled Spatial Transformation With Deterministic Mapping Function

1.1 Introduction

In Chapter 2, we elaborated on the details of the patient-adaptive ECG classification framework. An important part of this methodology is to design a personalized classifier with a deviation analysis module. The performance of the deviation analysis module depends on the geometry of different classes of data samples in the feature space. In Chapter 3, we proposed a novel spatial transformation method based on MOPSO to achieve a desired symmetry in the transformed feature space to boost the performance of the deviation analysis module. This methodology enables us to further process the normal samples and identify fuzzy states between the normal and abnormal states. However, the proposed method uses an iterative optimization which may have high computational complexity. In this chapter, we introduce another deterministic spatial transformation method to model the fuzzy state of samples with a more tractable and analytical solution.

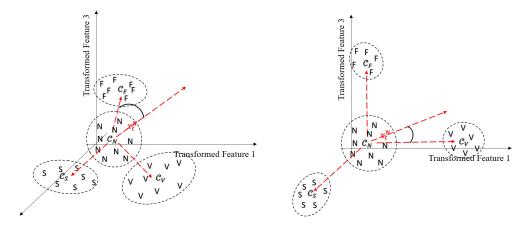


Figure 1.1: Left: illustration of the clustering topology in the original feature space; Right: illustration of the clustering topology in the feature space transformed with an optimized mapping function. Razi: it is not clear what you mean here by simple, is it unoptimized, why not just saying the original space before transmission!

While the aforementioned method in Chapter 3 yields the desired capacity of predicting upcoming abnormalities, the interpretation of the mechanisms of the system is not straightforward and easily tractable, thus it is not easily generalizable to a broader range of applications in the biomedical signal processing. The main objective of this chapter is to develop a deterministic transformation with a stronger prediction power based on a controlled spatial topology studied in Chapter 3. In this method, we further optimize the topology and spatial geometry of the clusters in the feature space by reducing the within-cluster variance after the spatial transformation. As shown in Fig. 1.1, through this improvement, the performance of the personalized classifier in terms of prediction accuracy can be further enhanced with the proposed method.

1.2 Hyper-Spherical Coordinates

In the previous chapter, the spatial transformation module is implemented with a polynomial kernel function and a heuristic optimization method. The system performance has been proven to be promising. In addition to the general drawbacks of heuristic method, it is not straightforward to select an appropriate base kernel as the core of spatial mapping function due to the high variety of kernel functions.

In order to address this issue, a novel deterministic spatial mapping function is proposed in this chapter based on *hyper-spherical* coordinates [REF: XXX (it is ok to repeat references, ans is ok to have web pages as REFs)]. Since hyper-spherical coordinates consist of angles and radii, these parameters in both original feature space and the desired target space are used to determine the mapping function.

The hyper-spherical coordinate system (n-dimensional spherical coordinate system) and its mapping to the *Cartesian* coordinate system are first introduced in [52]. If \mathbf{x} is a sample vector in a *n*-dimensional feature space, with its Cartesian coordinates (ξ_1 , ξ_2 ... ξ_n), then its corresponding hyper-spherical coordinates can be obtained through Eq.1.1, which is originally derived through its reverse mapping (Eq.1.2) using equation: $\sin(\arccos(x)) = \sqrt{1-x^2}$.

$$r = \sqrt{\xi_n^2 + \xi_{n-1}^2 + \dots + \xi_2^2 + \xi_1^2}$$

$$\theta_1 = \arccos \frac{\xi_1}{\sqrt{\xi_n^2 + \xi_{n-1}^2 + \dots + \xi_1^2}}$$

$$\theta_2 = \arccos \frac{\xi_2}{\sqrt{\xi_n^2 + \xi_{n-1}^2 + \dots + \xi_2^2}}$$

$$\vdots$$

$$\theta_{n-2} = \arccos \frac{\xi_{n-2}}{\sqrt{\xi_n^2 + \xi_{n-1}^2 + \xi_{n-2}^2}}$$

$$\theta_{n-1} = \begin{cases} \arccos \frac{\xi_{n-1}}{\sqrt{\xi_n^2 + \xi_{n-1}^2}} & \xi_n \ge 0 \\ -\arccos \frac{\xi_{n-1}}{\sqrt{\xi_n^2 + \xi_{n-1}^2}} & \xi_n < 0 \end{cases}$$

$$(1.1)$$

$$\xi_{1} = r \cos(\theta_{1})$$

$$\xi_{2} = r \sin(\theta_{1}) \cos(\theta_{2})$$

$$\xi_{3} = r \sin(\theta_{1}) \sin(\theta_{2}) \cos(\theta_{3})$$

$$\vdots$$

$$\xi_{n-1} = r \sin(\theta_{1}) \cdots \sin(\theta_{n-2}) \cos(\theta_{n-1})$$

$$\xi_{n} = r \sin(\theta_{1}) \cdots \sin(\theta_{n-2}) \sin(\theta_{n-1}). \tag{1.2}$$

where
$$0 \le \theta_j \le \pi$$
, $j = 1, ..., n - 2$; $0 \le \theta_{n-1} \le 2\pi$; $0 \le r < \infty$

1.3 Orthogonalization

To simplify the algorithm, in this chapter, the topology of clusters in the feature space is approximated by their centroid locations, denoted by $\mathbf{c}_N^k, \mathbf{c}_{\mathcal{V}}, \mathbf{c}_{\mathcal{S}}, \mathbf{c}_{\mathcal{F}}$, respectively for the normal cluster and the three abnormality clusters of type V, S, and F. Furthermore, as we assume that the samples with fuzzy states deviate from the normal cluster to a specific abnormality cluster, the spatial topology stays unchanged if the normal centroid is simply translated to the origin of the Cartesian coordinate system. The clustering topology in the original feature space can be equivalently represented by the following matrix with three row vectors:

$$C = \begin{bmatrix} \mathbf{c}_{\mathcal{V}} - \mathbf{c}_{N}^{k} \\ \mathbf{c}_{\mathcal{S}} - \mathbf{c}_{N}^{k} \\ \mathbf{c}_{\mathcal{F}} - \mathbf{c}_{N}^{k} \end{bmatrix} = \begin{bmatrix} \mathbf{v}_{\mathcal{V}\mathcal{N}} \\ \mathbf{v}_{\mathcal{S}\mathcal{N}} \\ \mathbf{v}_{\mathcal{F}\mathcal{N}} \end{bmatrix}$$
(1.3)

RZ: I used lower case v for vectors, which is more consistent with other definitions. Capital letter are more commonly used for matrices.

As shown in Fig. 1.1, in order to improve the performance of the personalized classifier and to avoid ambiguity in quantifying the deviations to different abnormality classes, a topology with a maximum separation between the three vectors of \mathcal{C} is preferred. To achieve a lower computational complexity in higher order dimensions, the algorithm aims at transforming vectors in \mathcal{C} to orthogonal vectors using a deterministic function. This approach not only simplifies the transformation derivations, it also ensures a full symmetry among abnormality clusters. As such, the problem reduces to finding a transformation that maps a set of vectors to a set of orthogonal vectors in the same space. Therefore, we can use the popular method of Gram-Schmidt orthogonalization method RZ: introduced or explained, choose the right term in [53,54]. Hence, in the first step of this stage, the three row vectors of \mathcal{C} representing the centroids of the three abnormality clusters are fed to the orthogonalization process as follows:

$$C^{\perp} = \operatorname{Gram-Schmidt}(C) = \begin{bmatrix} \mathbf{v}_{\mathcal{V}\mathcal{N}}^{\perp} \\ \mathbf{v}_{\mathcal{S}\mathcal{N}}^{\perp} \\ \mathbf{v}_{\mathcal{F}\mathcal{N}}^{\perp}, \end{bmatrix}$$
(1.4)

where \mathcal{C}^{\perp} is the matrix of orthogonalized vectors in the Cartesian coordinate system. The hyper-spherical coordinates \mathcal{C}^{\perp}_* of these orthogonalized vectors are calculated subsequently

using Eq.1.1. After this step, the orthogonalized vectors in the hyper-spherical coordinate system are obtained as formulated in Eq. 1.5. The j^{th} angular dimension denoted as θ_j^{\perp} and the radius is noted as r^{\perp} .

$$\mathcal{C}_{*}^{\perp} = \begin{bmatrix}
r_{\mathcal{V}\mathcal{N}}^{\perp} & \theta_{1_{\mathcal{V}\mathcal{N}}}^{\perp} & \dots & \theta_{n-1_{\mathcal{V}\mathcal{N}}}^{\perp} \\
r_{\mathcal{S}\mathcal{N}}^{\perp} & \theta_{1_{\mathcal{S}\mathcal{N}}}^{\perp} & \dots & \theta_{n-1_{\mathcal{S}\mathcal{N}}}^{\perp} \\
r_{\mathcal{F}\mathcal{N}}^{\perp} & \theta_{1_{\mathcal{F}\mathcal{N}}}^{\perp} & \dots & \theta_{n-1_{\mathcal{F}\mathcal{N}}}^{\perp}
\end{bmatrix}$$
(1.5)

1.4 Spatial Mapping Function

After obtaining the original hyper-spherical coordinates of $[\mathbf{v}_{\mathcal{V}\mathcal{N}}, \mathbf{v}_{\mathcal{S}\mathcal{N}}, \mathbf{v}_{\mathcal{F}\mathcal{N}}]^T$ and the orthogonalized hyper-spherical coordinates $[\mathbf{v}_{\mathcal{V}\mathcal{N}}^{\perp}, \mathbf{v}_{\mathcal{S}\mathcal{N}}^{\perp}, \mathbf{v}_{\mathcal{F}\mathcal{N}}^{\perp}]^T$, the goal is to design a mapping function $\mathbf{F}: \mathbf{R}^n \to \mathbf{R}^n$ from the original coordinates to the orthogonal ones which exhibit the desired clustering topology, such that Equi. 1.6 holds.

In the Gram-Schmidt algorithm, the very first vector serves as a reference vector and remains unchanged in the orthogonalization process, namely $\mathbf{v}_{\mathcal{V}\mathcal{N}} = \mathbf{v}_{\mathcal{V}\mathcal{N}}^{\perp}$. Therefore the mapping function \mathbf{F} shall satisfy the following equations:

$$\mathbf{F}(\mathbf{v}_{SN} - \mathbf{v}_{VN}) = \mathbf{v}_{SN}^{\perp} - \mathbf{v}_{VN}^{\perp} = \mathbf{v}_{SN}^{\perp} - \mathbf{v}_{VN}$$

$$\mathbf{F}(\mathbf{v}_{FN} - \mathbf{v}_{VN}) = \mathbf{v}_{FN}^{\perp} - \mathbf{v}_{VN}^{\perp} = \mathbf{v}_{FN}^{\perp} - \mathbf{v}_{VN}$$
(1.6)

Furthermore, since the orthogonality of vectors is independent to their radii r, we only need to design \mathbf{F} for the n-1 angular dimensions $(\theta_1, \ldots, \theta_{n-1})$, and the coordinate r remains unchanged after the mapping. Consequently, the function \mathbf{F} can be decomposed

into n-1 functions: $f_i: \mathbf{R} \to \mathbf{R}, i = 1...n-1$ with constraints in Eq.1.6 as well as two additional extreme points defining the range of input and output values. We use the notation $\mathbf{v}_{SN} - \mathbf{v}_{VN} = \Delta_{SV}$ and denote the i^{th} angular dimension of Δ_{SV} as $\delta_{i_{SV}}$. We use similar notations for the other dimensions, (e.g. $\mathbf{v}_{\mathcal{F}N} - \mathbf{v}_{VN}$). Hence for each angular dimension i, f_i is determined by $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{\mathcal{F}V}}, \delta_{i_{\mathcal{F}V}}^{\perp})$, along with the two extreme boundary points defining the domain and range of the functions.

In order to maintain the simplicity and the linearity of the mapping function, the functions f_i should be continuous and monotonic. For this purpose, the valid range of angular dimensions (after considering the periodicity of these functions) is used to determine the boundary constrains; therefore the problem is transformed into a curve fitting problem.

RZ: the text was not clear, it now reads much better. For example, if linking $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$ results in a monotonically increasing function, noting the fact that the range of n-2 first angular dimensions is the interval $[0,\pi]$, then the two extreme boundary points would be (0,0) and (π,π) , as depicted in Fig. 1.2. Conversely, for a monotonically increasing function, the extreme boundary points would be $(\pi,0)$ and $(0i,\pi)$. These rules apply for the functions defined for the n-2 first angular dimensions. Similar rules apply to the last angular dimension with only one consideration that the period is 2π instead of π . Since $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$ represent the orthogonalization process, they are decisive for this process. So we call these two points as well as the two extreme boundary points as target points of the angular dimensions in the following sections.

The simplest candidate function for f_i , which connects the two boundary points and the two target points in the 2-D plane would be a linear spline as shown in Fig.1.2.

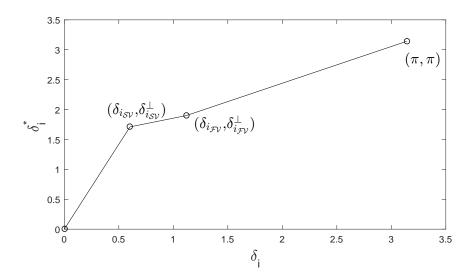


Figure 1.2: The simple mapping function for one angular dimension which maps the target points in the original space to the desired target points. Target points include the two extreme boundary points (0,0), and π,π) as well as (XXX,XXX) and (XXX,XXX) to yield the desired mapping. add (0,0) and the label target and extreme boundary points to the figure.

1.5 Optimized Mapping Function

The mapping function in Fig.1.2 exhibits the two desired properties of monotonicity and continuity for an ideal mapping function f_i . However, this simple peicewise linear mapping function suffers from some drawbacks. Firstlt, the function is not differentiable at the target points $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$, which can lead to severe cluster deformation. Secondly, the mapping function is applied to the angular dimensions, the discontinuity of the derivative of the function can cause the convex-clusters to be mapped into non-convex clusters in the Cartesian coordinate system. In order to avoid deformations and to preserve maximal similarity with the original clustering geometry, it is desired to use a more linear functions with continuous derivatives. Also, in order to provide maximal separation among clusters and to concentrate clusters into disjoint non-overlapping clusters, it is beneficial to map the regions between two target points into a region with equal or even smaller range in

the Cartesian coordinate system. In other words, it is desired to keep the samples close to the centroids in the mapped space as much as possible. This property is achieved by using nonlinear functions. However, there is a trade-off between the level of concentration and the linearity, which needs to be carefully addressed. In order to avoid deformation while providing maximal separation, an optimized mapping function is proposed in this section.

In order to accommodate satisfy the above-mentioned requirements, it is desired to find a function, which satisfies the following mathematical conditions:

- the function is differentiable everywhere (continuous first derivative);
- the derivative of the function is small at the target points, which correspond to the centroids of clusters in the original and transformed space, i.e. $\delta_{i_{SV}}$ and $\delta_{i_{FV}}$ in Fig. XXX;
- the derivative of the function is large at the boundaries of two regions (point $(\epsilon_{i_{sv}}, XXX)$) in Fig. XX.1.3;

Therefore, we propose to use the basis function p, which satisfies the aforementioned conditions. Why you call it basis function. Basis function has a different meaning in linear analysis. I suggest you use the term basic function with adjustable parameters. Each target point is associated with a basic function. The function p is composed of two constitutive functions: h(x) and g(x), respectively defined in two regions: i) from the target point to the upper boundary point, ii) from the lower boundary point to the next target point. [RZ: show h(x) and g(x) in Fig. 1.3.]. Before defining the two functions h(x) and g(x), we need to define the boundaries between two consecutive target points $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$, in which these functions are defined (i.g. $\epsilon_X X X X$). We simply choose the midpoint as the boundary points. For instance, we have $\epsilon = (\delta_1 + \delta_2)/2$ [RZ:Write the correct equation] as

stated in Equi. 1.7. The lower and upper boundary points are denoted respectively as (γ, γ^{\perp}) and $(\epsilon, \epsilon^{\perp})$. To ensure the continuity of the mapping function f, the boundaries and target points should satisfy:

$$(\epsilon_{i_{\mathcal{S}\mathcal{V}}}, \epsilon_{i_{\mathcal{S}\mathcal{V}}}^{\perp}) = (\gamma_{i_{\mathcal{F}\mathcal{V}}}, \gamma_{i_{\mathcal{F}\mathcal{V}}}^{\perp}) = (\frac{\delta_{i_{\mathcal{S}\mathcal{V}}} + \delta_{i_{\mathcal{F}\mathcal{V}}}}{2}, \frac{\delta_{i_{\mathcal{S}\mathcal{V}}}^{\perp} + \delta_{i_{\mathcal{F}\mathcal{V}}}^{\perp}}{2})$$
(1.7)

The two piece-wise functions h(x) and g(x) are defined as use the correct name: the inverse of logit/sigmoid function as follows:

$$K_h = \frac{\epsilon^{\perp} - \delta^{\perp}}{e^{\alpha(\epsilon - \delta, 0)^{+}} - 1}$$

$$h(x) = K_h[e^{\alpha(x - \delta, 0)^{+}} - 1] + \delta^{\perp}$$
(1.8)

$$K_g = \frac{\gamma^{\perp} - \delta^{\perp}}{e^{\alpha(-\gamma + \delta, 0)^{+}} - 1}$$

$$g(x) = K_g[e^{\alpha(\delta - x, 0)^{+}} - 1] + \delta^{\perp}$$

$$(1.9)$$

RZ: revised up to here 2.

To demonstrate this process, we first apply p on two target points $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$ and $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$. This will result in a smooth curve as shown in Fig.1.3

Therefore, if we apply the piecewise interpolate function p at all four target points: $(\delta_{i_{SV}}, \delta_{i_{SV}}^{\perp})$, $(\delta_{i_{FV}}, \delta_{i_{FV}}^{\perp})$ and two boundary points of angular dimension, the final result is shown in Fig.1.4 In the training process, abnormal data from DS1 and personalized normal cluster are used to

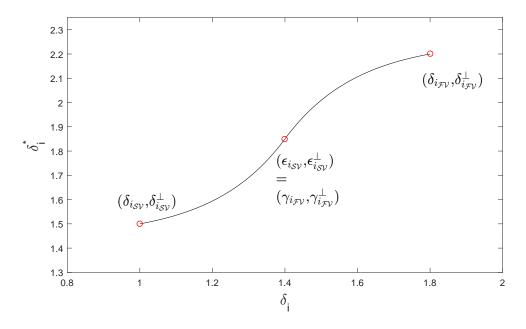


Figure 1.3: Optimized Piece-wise Interpolated Function p. I understand this figure, but is also misleading since the target points are at two ends and the boundary is in middle (the opposite is better). To solve this issue you can plot two adjacent regions and use tags and labels to show the target points δ and edge nodes/boundary nodes ϵ .

determine the mapping function. In the predicting process, the hyper-spherical coordinate of a new ECG sample is calculated and the mapping function is applied on its hyper-spherical coordinate. After this step, we calculate the Cartesian coordinate of the transformed data. This final result in Cartesian coordinate is then fed into Eq. ?? to generate the corresponding type of yellow alarm.

1.6 Experimental Results

In this section, the performance of proposed method is evaluated according to two aspects. We will first analyze the classification performance of the system and demonstrate the comparison with other representative ECG classifiers. Furthermore, the classification results are partitioned into two sets: red alarms generated by global classifier and final results by

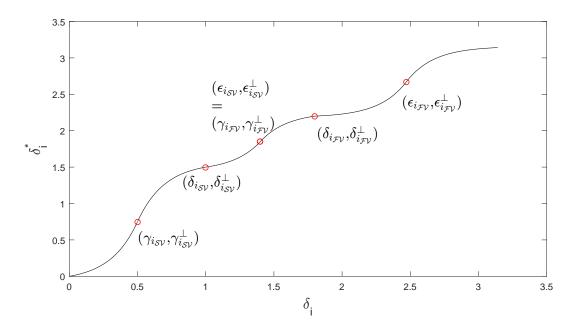


Figure 1.4: Optimized Mapping Function f.

combining yellow and red alarms. In this way, the impacts of personalized classifier on final results can be revealed. Finally, the prediction performance which is representative in the proposed system is evaluated.

1.6.1 Classification Performance

The experimental results are evaluated with the classification performance of 4 AAMI ECG classes using the test subset of MITBIH Arrhythmia DS2. Originally DS2 contains 15357 samples after feature extraction. While training Personal Classifier, the first 20% of total normal samples serve as initialization samples for personalized dynamic normal cluster. Therefore, all samples before the last initialization normal sample should be excluded from test set for each records. Consequently, the actual test set contains 12414 samples in total consisting of 10105 type-N, 1702 type-V, 508 type-S and 99 type-F samples.

To present the result, we select weighted k-Nearest Neighbors where k=10 as global classifier

Table 1.1: Cumulated Confusion Matrix for All Records in DS2

	Ground Truth										
		N	V	S	F						
Result	N	9255(10076)	21(38)	72(90)	1(5)						
	V	657(22)	1678(1663)	8(2)	9(7)						
	S	71(6)	3(1)	417(416)	0(0)						
	F	122(1)	0(0)	11(0)	89(87)						

because it's comparatively simple and representative among low complexity models. The parameter α in the deviation detection module is set to 1 for test purpose.

Table 1.1 summarized the cumulated confusion matrix for all records in the test set. In order to compare the result of global classifier and combined result, the sample numbers are presented in the following format: combined(globaled). In order to measure classification performance, we adopted three metrics proposed in [12,13,16]: accuracy(Ac), sensitivity(Se), specificity(Sp). All three metrics are calculated based on true positive TP, false positive FP, false negative FN and true negative TN in a binary confusion matrix. Therefore all four metrics are calculated for each class by converting the 4x4 matrix to a 2x2 matrix.

While cumulated classification results are demonstrated in Table 1.1, the robustness of the proposed method should be evaluated based on the performance variation over 22 test records in DS2. Hence medians and IQRs(interquartile range) for each metric and each class are included in Table 1.2 to represent the robustness of proposed methods. The lower variation between performances measured on different tapes, the more robust the system is. In Table 1.2, we observe that among all abnormal classes, the proposed method demonstrates stable performance on class V but less stable on class S and F.

As MITDB is widely used to verify ECG classifier performance, we compared the proposed system with five significant methods proposed in literature. According to AAMI standards, ECG classifier performance should be evaluated over the binary classification performance of Ventricular(V) versus non-V types and Supraventricular(S) versus non-S types. For methods

ods proposed in literature, same evaluation metrics are deployed on records from MITDB. To standardize the metrics, we select 11 common ECG records from all 5 methods and compared the median of each classification metrics over these 11 records. The comparison results are demonstrated in Table 1.3. Generally speaking, the proposed method shows higher sensitivity for both V type and S type. Especially for S type, the proposed method has advantage on all three metrics over other five methods in the literature.

1.6.2 Prediction Performance

As an important characteristic of the proposed methods, yellow alarms triggered by personalized classifier after feature space reshaping indicate a higher probability of observing subsequent abnormalities. In order to verify this functionality, all beats following a yellow alarm is investigated for each yellow alarm. Abnormality type which occurs the earliest within the window is recorded. Similar to confusion matrix for classification evaluation, the performance of prediction can be summarized by a confusion matrix with the 3 abnormal types. Probabilities of observing a certain type of abnormal beats after a yellow alarm is calculated using the prediction confusion matrix and compared to the prior probability of observing the same type of abnormality. This process is formulated in the following two equations:

Table 1.2: Classification Performance and Within-Set Variation of Proposed System

statistics	N			V			S			F		
Statistics	Ac	Se	Sp									
cumulated	92.4	91.59	95.93	94.38	98.59	93.71	98.67	82.09	99.38	98.85	89.9	98.92
median	94.45	92.21	95.42	96.17	99.55	95.71	99.38	80.65	99.84	99.11	90.91	99.11
IQR	6.33	10.08	11.91	5.17	1.64	8.62	1.76	19.35	0.61	1.58	23.33	1.49

Table 1.3: V and S classification performance compared with five algorithms in literature using 11 common records in MITDB

Methods		VEB		SVEB			
Methods	Ac	Se	Sp	Ac	Se	Sp	
Proposed	96.6	98.2	92.4	98.63	88.89	99.41	
Hu et al. [12]	94.8	78.9	96.8	N/A	N/A	N/A	
de Chazal et al. [9]	96.4	77.5	N/A	N/A	N/A	N/A	
Jiang and Kong [15]	98.8	78.9	96.8	97.5	74.9	98.8	
Ince <i>et al.</i> [16]	97.9	90.3	98.8	96.1	81.8	98.5	
Kiranyaz et al. [17]	98.9	95.9	99.4	96.4	68.8	99.5	

Table 1.4: predictive probability versus prior probability without windowing

	# of	predi	cted ground truth	% of predicted ground truth			
	V	S	F	V	S	F	
yellow alarm	V	467	122	14	77.45	20.23	2.32
	S	36	15	0	70.59	28.41	0
	F	40	60	5	38.10	57.14	4.76
total		543	197	19	71.54	25.96	2.50

$$P(\hat{y}_{k+i} = X_r | \hat{y}_k = X_y) = \frac{\text{# of } y_{k+i} = X \text{ after } \hat{y}_k = X_y}{\text{# of true alarms after } \hat{y}_k = X_y}$$

$$P(\hat{y}_{k+i} = X_r) = \frac{\text{# of true alarm of type } X (y_k = X)}{\text{# of all true alarms}}$$
(1.10)

The power of predicting each type of abnormalities is evaluated by comparing $P(\hat{y}_{k+i} = X_r | \hat{y}_k = X_y)$ and $P(\hat{y}_{k+i} = X_r)$. A shown in Table 1.4, the probability of observing a certain type of abnormalities after a yellow alarm is higher than its prior for each type of abnormalities. For example, without knowing the type of a yellow alarm, the probability of observing a type V sample is 71.54% while the probability of observing a type V sample given a type V yellow alarm was triggered is 77.45%. The improvement are consistent among all three types of abnormalities but the system has stronger predicting capacity for type S.

Table 1.5: predictive probability versus prior probability within 10 beats' window

	# of	predi	cted ground truth	% of predicted ground truth			
		V	S	F	V	S	F
	V	290	85	12	74.94	21.96	3.10
yellow alarm	S	22	13	0	62.86	37.14	0
	F	29	37	6	40.28	51.39	8.33
total		341	135	18	69.03	27.32	3.64

In order to study the time window in which real abnormality occurs after yellow alarms, we also studied a window of 10 consecutive samples following a yellow alarm. Similarly, prior and posterior probabilities are compared to evaluate the performance as shown in Table.1.5.

Compared with the result without windowing, the predicting performance within 10 beats window shows that the proposed algorithm can better predict the occurrence of abnormalities in a certain time window. Especially for type S, the probability of observing a type S sample within 10 beats after a yellow alarm is 27.32% while given that the yellow alarm is type S, the posterior probability is raise to 37.14%. With almost 10% increase, it's proved that the yellow alarm types are informative. The results shows that same improvements are made within the 10-sample window as well. In general, the predicting performance are promising, indicating the efficiency of personalized classifier and deviation analysis.

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