

NEW ADAPTIVE ALGORITHMS FOR OPTIMAL FEATURE EXTRACTION FROM GAUSSIAN DATA

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Abstract: In this paper, we present new adaptive learning algorithms to extract optimal features from multidimensional Gaussian data while preserving class separability. For this purpose, we introduce new adaptive algorithms for the computation of the square root of the inverse covariance matrix $\Sigma^{-1/2}$. We prove the convergence of the adaptive algorithms by introducing the related cost function and discussing about its properties and initial conditions. Adaptive nature of the new feature extraction method makes it appropriate for on-line signal processing and pattern recognition applications. Experimental results using two-class multidimensional Gaussian data demonstrated the effectiveness of the new adaptive feature extraction method.

1 INTRODUCTION

Feature extraction is generally considered as a process of mapping the original measurements into a more effective feature space. When we have two or more classes, feature extraction consists of choosing those features which are the most effective for preserving class separability in addition to dimension reduction (Theodoridis, 2003). One of the most used techniques for this purpose is linear discriminant analysis (LDA) algorithm. LDA algorithm has been widely used in signal processing and pattern recognition applications in which feature extraction is inevitable, such as face and gesture recognition and hyper-spectral image processing (Chang and Ren, 2000; Chen *et al.*, 2000; Lu *et al.* 2003) Conventional LDA algorithm is used only in off-line applications. However, the needs for dimensionality reduction in real time applications such as on-line face recognition, motivated researchers to introduce adaptive versions of LDA. Chatterjee and Roychowdhury presented an adaptive algorithm and a self-organizing LDA network for feature extraction from Gaussian data (Chatterjee and Roychowdhury, 1997). They introduced an adaptive method for computation of $\Sigma^{-1/2}$ in which Σ is the symmetric positive definite scattering matrix of a random vector sequence. However, they didn't introduce any cost function related to their

adaptive algorithm. Therefore, they used the stochastic approximation theory in order to prove the convergence of their adaptive equation and outlined networks for feature extraction. On the other hand, the approach presented in (Chatterjee and Roychowdhury, 1997) suffers from low convergence rate. Recently, Abrishami Moghaddam *et al.* (2003; 2005) proposed three new adaptive methods based on steepest descent, conjugate direction and Newton-Raphson optimization techniques to hasten convergence of the adaptive algorithm.

In this study, we present new adaptive algorithms for the computation of $\Sigma^{-1/2}$. Furthermore, we introduce a cost function related to these algorithms and prove their convergence by discussing about the properties and initial conditions of this cost function. Existence of the cost function and its differentiability facilitate the convergence analysis of the new adaptive algorithms without using complicated stochastic approximation theory. We will show effectiveness of these new adaptive algorithms for extracting optimal features from two-class multi-dimensional Gaussian sequences.

The organization of the paper is as follows. The next section describes the fundamentals of optimal feature extraction from Gaussian data. Section 3, presents the new adaptive equations and analyzes their convergence by introducing the related cost function and discussing about initial conditions. Section 4, is devoted to simulations and

experimental results. Finally, concluding remarks are given in section 5.

2 OPTIMAL FEATURES FOR GAUSSIAN DATA

Let $\{\omega_1, \omega_2, \dots, \omega_L\}$ be the L classes in which our patterns belong and $\mathbf{x} \in \mathfrak{R}^n$ be a pattern vector whose mixture distribution is given by $p(\mathbf{x})$. In a sequel, it is assumed that a priori probabilities $P(\omega_i)$, $i=1, \dots, L$, are known. If they are not explicitly known, they can easily be estimated from the available training vectors. For example, if N is the total number of available training patterns and N_i ($i=1, \dots, L$) of them belong to ω_i , then $P(\omega_i) \approx N_i/N$. Consider conditional probability densities $p(\mathbf{x} | \omega_i)$, $i=1, \dots, L$ and posterior probabilities $P(\omega_i | \mathbf{x})$, $i=1, \dots, L$ are known. Using Bayes classification rule, we can state that the pattern \mathbf{x} is classified to ω_i if $P(\omega_i | \mathbf{x}) > P(\omega_j | \mathbf{x})$, $i, j=1, \dots, L$ and $j \neq i$

In other words, the L a posteriori probability functions, mentioned above, are sufficient statistics, and carry all information for classification in the Bayes sense. The Bayes classifier in this feature space is a piecewise bisector classifier which is its simplest form (Fukunaga, 1990). Gaussian distribution in general has a density function in the following form

$$N(\mathbf{m}, \Sigma) = \frac{1}{(2\pi)^{\frac{n}{2}} |\Sigma|^{\frac{1}{2}}} e^{-\frac{1}{2}d^2(\mathbf{x})} \quad (1)$$

where the distance function $d^2(\mathbf{x})$, is defined by:

$$d^2(\mathbf{x}) = (\mathbf{x} - \mathbf{m})^t \Sigma^{-1} (\mathbf{x} - \mathbf{m}) \quad (2)$$

Which $\mathbf{x} \in \mathfrak{R}^n$ is a random vector, Σ is a $n \times n$ symmetric covariance matrix and \mathbf{m} is a $n \times 1$ vector denoted mean value of the random sequence. Considering the feature:

$$\ln P(\omega_i | \mathbf{x}) = \ln p(\mathbf{x} | \omega_i) + \ln P(\omega_i) - \ln p(\mathbf{x}) \quad (3)$$

for class ω_i , $i=1, \dots, L$, it will appear that, $\ln p(\mathbf{x} | \omega_i)$ is the relevant feature for class ω_i (recalling that in feature extraction, additive and multiplicative constants do not modify the subspace onto which the distributions are mapped). Supposing unimodal Gaussian distribution, the feature

In $p(\mathbf{x} | \omega_i)$ reduces to a quadratic function $f_i(\mathbf{x})$, defined as:

$$f_i(\mathbf{x}) = (\mathbf{x} - \mathbf{m}_i)^t \Sigma_i^{-1} (\mathbf{x} - \mathbf{m}_i), \quad i=1, \dots, L \quad (4)$$

where \mathbf{m}_i and Σ_i are the class ω_i 's mean value and covariance matrix, respectively. The function $f_i(\mathbf{x})$, can be expressed in the form of a norm function (Fukunaga, 1990):

$$f_i(\mathbf{x}) = \|\Sigma_i^{-\frac{1}{2}} (\mathbf{x} - \mathbf{m}_i)\|^2, \quad i=1, \dots, L \quad (5)$$

From the above discussion, it is clear that function $f_i(\mathbf{x})$ is the sufficient information for classification of Gaussian data with minimum Bayes error. In other words, after computation of $f_i(\mathbf{x})$ for $i=1, \dots, L$, it is easy to decide about classification of the unknown vector \mathbf{x} . Generally speaking, in on-line applications, the values of $\Sigma^{-1/2}$ and \mathbf{m}_i are unknown. Therefore, we should find a rule for adaptive estimation of these values and compute $f_i(\mathbf{x})$. In the next section, a new method for adaptive computation of $\Sigma^{-1/2}$ will be presented. In addition, we introduce a cost function related to this adaptive equation, and use it for proving its convergence.

3 ADAPTIVE COMPUTATION OF $\Sigma^{-1/2}$ AND CONVERGENCE PROOF

We define the cost function $J(\mathbf{W})$ with parameter \mathbf{W} $J: \mathfrak{R}^{n \times n} \rightarrow \mathfrak{R}$ as follows:

$$J(\mathbf{W}) = \frac{\text{tr}(\mathbf{W}^3 \mathbf{X} \mathbf{X}^t)}{3} - \text{tr}(\mathbf{W}) \quad (6)$$

$J(\mathbf{W})$, is a continuous function with respect to \mathbf{W} . The expected value of J (for constant \mathbf{W}) is given by:

$$E(J(\mathbf{W})) = \frac{\text{tr}(\mathbf{W}^3 \Sigma)}{3} - \text{tr}(\mathbf{W}) \quad (7)$$

where Σ is the covariance matrix. The first derivative of $E(J)$ is computed as follows (assuming that \mathbf{W} is a symmetric matrix) (Magnus, Neudecker, 1999):

$$\frac{\partial E(J(\mathbf{W}))}{\partial \mathbf{W}} = (\mathbf{W}^2 \Sigma + \mathbf{W} \Sigma \mathbf{W} + \Sigma \mathbf{W}^2) / 3 - \mathbf{I} \quad (8)$$

The unique zero solution of (8) is $\Sigma^{-1/2}$, the second derivative of the expected value of the cost function is equal to (Magnus, Neudecker, 1999):

$$\frac{\partial^2 E(J(\mathbf{W}))}{\partial^2 \mathbf{W}} = 2(\mathbf{I} \otimes \Sigma \mathbf{W}) + 2(\Sigma \mathbf{W} \otimes \mathbf{I}) + \mathbf{W} \otimes \Sigma + \Sigma \otimes \mathbf{W} \quad (9)$$

In (9) it is assumed that \mathbf{w} is a symmetric matrix that it commutes with Σ . If we substitute \mathbf{w} in (9) with $\Sigma^{-1/2}$, the answer will be:

$$\left. \frac{\partial^2 E(J(\mathbf{W}))}{\partial^2 \mathbf{W}} \right|_{\mathbf{w}=\Sigma^{-1/2}} = 2(\mathbf{I} \otimes \Sigma^{1/2}) + 2(\Sigma^{1/2} \otimes \mathbf{I}) + \Sigma^{-1/2} \otimes \Sigma + \Sigma \otimes \Sigma^{-1/2} \quad (10)$$

where (10) is a positive definite matrix. From the above discussion, it can be concluded that the cost function $J(\mathbf{w})$ has a minimum that occurs at $\Sigma^{-1/2}$ (Magnus, Neudecker, 1999). Using the gradient descent optimization method (Widrow, Stearns, 1985; Hagan, Demuth, 2002), we obtain the following adaptive equation for the computation of $\Sigma^{-1/2}$:

$$\mathbf{W}_{k+1} = \mathbf{W}_k + \eta_k \left(-\frac{\partial J(\mathbf{W})}{\partial \mathbf{W}} \right) = \mathbf{W}_k + \eta_k (I - (\mathbf{W}_k^2 \mathbf{x}_{k+1} \mathbf{x}_{k+1}' + \mathbf{W}_k \mathbf{x}_{k+1} \mathbf{x}_{k+1}' \mathbf{W}_k + \mathbf{x}_{k+1} \mathbf{x}_{k+1}' \mathbf{W}_k^2) / 3) \quad (11)$$

In (11), \mathbf{W}_{k+1} is the estimation of $\Sigma^{-1/2}$ in $k+1$ -th iteration. η_k is the step size and \mathbf{x}_{k+1} is the input vector at iteration $k+1$. Equation (11) updates the estimation of $\Sigma^{-1/2}$, using the last estimation and the present input vector. The only constraint on (11) is its initial condition. That means \mathbf{W}_0 must be a symmetric and positive definite matrix satisfying $\mathbf{W}_0 \Sigma = \Sigma \mathbf{W}_0$. It is quite easy to prove that if \mathbf{W}_0 is a symmetric and positive definite matrix, then all values of \mathbf{W}_i ($i=2,3,\dots$) will be symmetric and positive definite. Therefore, the final estimation also will have these properties (which are essential for covariance matrix). To avoid confusion for choosing the initial value \mathbf{W}_0 , we considered \mathbf{W}_0 equal to identity matrix multiplied by a positive constant ($\mathbf{W}_0 = \alpha \mathbf{I}$).

According to the result reported by (Kushner, Clark, 1978; Benveniste, Metivier, 1990), the stochastic gradient algorithm in the form of:

$$\mathbf{W}_{k+1} = \mathbf{W}_k - \eta_k f(\theta_k, Y_{k+1}) \quad (12)$$

where $f(\theta, y) = \text{grad}_{\theta} F(\theta, y)$ and $(Y_k)_{k>0}$ are independent identically distributed \mathfrak{R}^n -valued random variables, converges almost surely towards a

solution of the minimization problem: $\min_{\theta} E(F(\theta, Y))$. As indicated in (12), in order to minimize $E(F(\theta, Y))$, the stochastic gradient algorithm uses the random variable $f(\theta, Y)$ instead of its expectation in the ordinary gradient method. The above argument is another approach for proving the convergence of (11) towards $\Sigma^{-1/2}$.

It is easy to show that if \mathbf{W}_0 considered a symmetric and positive definite matrix which satisfy $\mathbf{W}_0 \Sigma = \Sigma \mathbf{W}_0$ then expected value of (11) will be equal to the following equations:

$$\begin{aligned} E(\mathbf{W}_{k+1}) &= \mathbf{W}_k + \eta_k (\mathbf{I} - \mathbf{W}_k^2 \Sigma) \\ &= \mathbf{W}_k + \eta_k (\mathbf{I} - \mathbf{W}_k \Sigma \mathbf{W}_k) \\ &= \mathbf{W}_k + \eta_k (\mathbf{I} - \Sigma \mathbf{W}_k^2) \end{aligned} \quad (13)$$

Therefore (11) is simplified to three more efficient forms as follows:

$$\mathbf{W}_{k+1} = \mathbf{W}_k + \eta_k (\mathbf{I} - \mathbf{W}_k^2 \mathbf{x}_{k+1} \mathbf{x}_{k+1}') \quad (14)$$

$$\mathbf{W}_{k+1} = \mathbf{W}_k + \eta_k (\mathbf{I} - \mathbf{W}_k^2 \mathbf{x}_{k+1} \mathbf{x}_{k+1}' \mathbf{W}_k) \quad (15)$$

$$\mathbf{W}_{k+1} = \mathbf{W}_k + \eta_k (\mathbf{I} - \mathbf{W}_k \mathbf{x}_{k+1} \mathbf{x}_{k+1}' \mathbf{W}_k) \quad (16)$$

Existence of the cost function for the new adaptive $\Sigma^{-1/2}$ algorithms has the following advantages compared to the former one (Chatterjee and Roychowdhury, 1997): *i*) it simplifies the task for proving the convergence; *ii*) it helps to evaluate the accuracy of the current solutions. For example, in the cases of different initial conditions and various learning rates, one is enable to evaluate which initial condition and learning rate outperform others. Furthermore, the former adaptive equation in (Chatterjee and Roychowdhury, 1997) uses a fix or monotonically decreasing learning rate which results in low convergence speed, but introducing a cost function related to the adaptive algorithm make it possible to determine the learning rate efficiently in every stage in order to increase the convergence rate.

There are different methods for adaptive estimation of the mean vector. The following equation was used in (Chatterjee and Roychowdhury, 1997; Abrishami Moghaddam *et al.*, 2003; 2005):

$$\mathbf{m}_{k+1} = \mathbf{m}_k + \eta_k (\mathbf{x}_{k+1} - \mathbf{m}_k) \quad (17)$$

where η_k satisfies Ljung assumptions (Ljung, 1977) for the step size. Alternatively, one may use the following equation (Ozawa *et al.*, 2005; Pang *et al.*, 2005):

$$\mathbf{m}_{k+1} = \frac{k}{k+1} \mathbf{m}_k + \frac{\mathbf{x}_{k+1}}{k+1} \quad (18)$$

For the experiments reported in the next section, we used (17) in order to estimate the mean value in each iteration.

4 SIMULATION RESULTS

In this section, we used on of the (14-16) described in the previous section to estimate $\Sigma^{-1/2}$ and extracted features from Gaussian data for classification.

4.1 Experiments on $\Sigma^{-1/2}$ Algorithm

In the first experiment, we compared the convergence of the new adaptive $\Sigma^{-1/2}$ algorithm with the algorithm proposed in (Chatterjee and Roychowdhury, 1997). We used the first covariance matrix in (Okada and Tomita, 1985), which is a 10×10 covariance matrix and multiplied it, by 20 (Figure 1). The ten eigenvalues of this matrix in descending order are 117.996, 55.644, 34.175, 7.873, 5.878, 1.743, 1.423, 1.213 and 1.007. Figure 2 compares the error of each algorithm as a function of sample number. As illustrated, the new algorithm can converges with an accelerated rate than the previous algorithm. We also compared the convergence rate of the new adaptive $\Sigma^{-1/2}$ algorithm in 4, 6, 8 and 10 dimensional spaces.

$$\Sigma=20 \begin{bmatrix} 0.091 & & & & & & & & & \\ 0.038 & 0.373 & & & & & & & & \\ -0.053 & 0.018 & 1.430 & & & & & & & \\ -0.005 & -0.028 & 0.017 & 0.084 & & & & & & \\ 0.010 & -0.011 & 0.055 & -0.005 & 0.071 & & & & & \\ -0.136 & -0.367 & -0.450 & 0.016 & 0.088 & 5.720 & & & & \\ 0.155 & 0.154 & -0.038 & 0.042 & 0.058 & -0.544 & 2.750 & & & \\ 0.030 & -0.057 & -0.298 & -0.022 & -0.069 & -0.248 & -0.343 & 1.450 & & \\ 0.002 & -0.031 & -0.041 & 0.001 & -0.008 & 0.005 & -0.011 & 0.078 & 0.067 & \\ 0.032 & -0.065 & -0.030 & 0.003 & 0.003 & 0.095 & -0.120 & 0.028 & 0.015 & 0.341 \end{bmatrix}$$

Figure 1: Sample covariance matrix used in $\Sigma^{-1/2}$ experiments.

We used the same covariance matrix as in the first experiment for generating 10 dimensional data and three other matrices were selected as the principal minors of that matrix.

In all experiments, we chose the initial value \mathbf{W}_0 equal to identity matrix multiplied by 0.6, and then using a sequence of Gaussian input data (training data) estimated $\Sigma^{-1/2}$. For each covariance matrix,

we generated 500 samples of zero-mean Gaussian data and estimated the $\Sigma^{-1/2}$ matrix using (14).

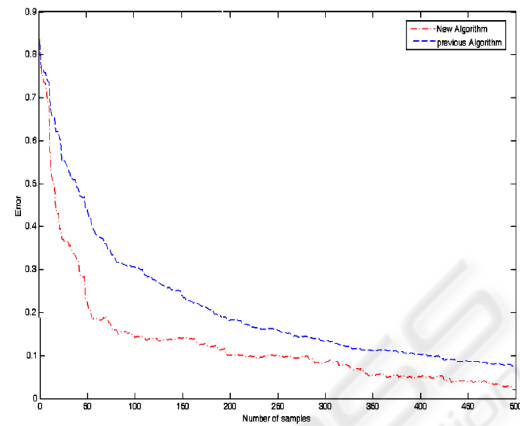


Figure 2: Comparison of convergence rate between new algorithm and previous algorithm.

For each experiment, at k -th iteration, we computed the error $e(k)$ between the estimated and actual $\Sigma^{-1/2}$ matrices by:

$$e(k) = \sqrt{\sum_{i=1}^n \sum_{j=1}^n (\mathbf{W}_{ij}(k) - \Sigma^{-1/2})^2} \quad (19)$$

For each covariance matrix, we computed the norm of error in every iteration. Figure 3 shows values of the error during iterations for each covariance matrix. The final values of error after 500 samples are error=0.169 for $d=10$, error=0.118 for $d=8$, error=0.102 for $d=6$ and error=0.0705 for $d=4$. As expected, the simulation results confirmed the convergence of (14) toward the $\Sigma^{-1/2}$. We repeated the same experiment using (15) and (16) and obtained similar results.

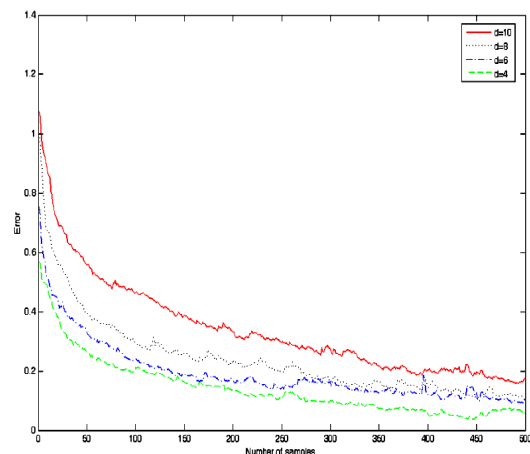


Figure 3: Convergence of the $\Sigma^{-1/2}$ algorithm for different covariance matrices.

4.2 Extracting Optimal Features from Two Class Gaussian Data

As discussed in section 2, it is apparent that for Gaussian data, the feature $f_i(\mathbf{x})$ is equal to $\|\Sigma_i^{-1/2}(\mathbf{x} - \mathbf{m}_i)\|^2$, using (14-16), (17) and the training sample sequence, we estimated $\Sigma^{-1/2}$ and \mathbf{m}_i .

For each training data belong to ω_i , we updated first $\Sigma^{-1/2}$ using (14-16) and then refreshed \mathbf{m}_i by applying (17), finally, we computed the norm of $\Sigma_i^{-1/2}(\mathbf{x}_i^k - \mathbf{m}_i)$. After computation of the mean value and $\Sigma^{-1/2}$ according to (5), it is possible to classify the next coming Gaussian data. At the end of this process, we compute the number of misclassifications. For testing the effectiveness of (14-16) in the case of two class Gaussian data, we generated 500 samples of 2 dimensional Gaussian data; each sample belonged to one of two classes with different covariance matrices and mean vectors. For each pattern \mathbf{X} , we extracted features $f_1(\mathbf{x})$ and $f_2(\mathbf{x})$. According to these optimal features, we transformed incoming Gaussian data into optimal feature space. Figures 4 and 5 show the comparison between samples in the original space and transformed samples in the optimal feature space. Two Gaussian classes ω_1 and ω_2 had the following parameters:

$$m_1 = \begin{bmatrix} -2 \\ 2 \end{bmatrix}, Q_1 = \begin{bmatrix} 3 & 2 \\ 2 & 3 \end{bmatrix}, m_2 = \begin{bmatrix} 2 \\ -2 \end{bmatrix}, Q_2 = \begin{bmatrix} 2 & 0 \\ 0 & 2 \end{bmatrix}$$

Figure 4 shows the distribution of samples from two classes. It is obvious that two classes are not linearly distinguishable. After estimation of $\Sigma^{-1/2}$ by (14-16) and estimation of \mathbf{m}_i by (17), we are able to extract f_1 and f_2 from the training data.

Figure 5 shows the transformed data in optimal feature space. It is apparent from Figure 5 that two Gaussian classes are linearly separable in the optimal feature space. In other words, in the optimal feature space, we can draw a straight line to separate two classes. However, in their original space; two classes are overlapped and are not linearly separable. By extracting optimal features, only 9 sample data among 1000 total sample, were misclassified by a linear classifier.

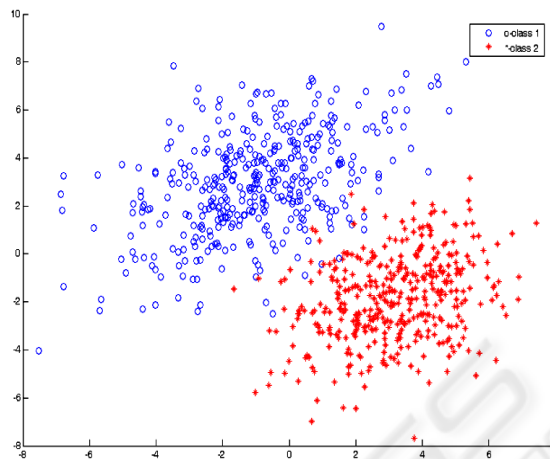


Figure 4: Distribution of two class Gaussian data in the original space.

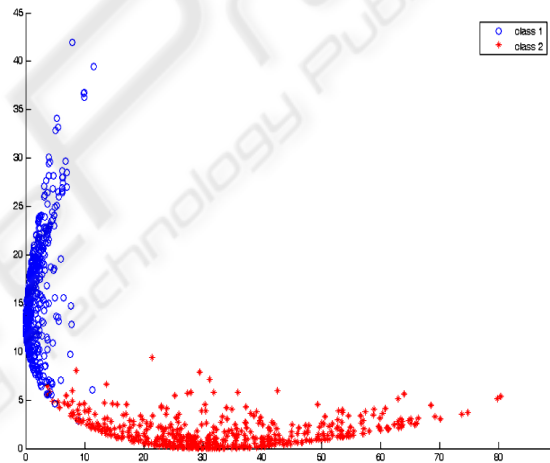


Figure 5: Distribution of two class Gaussian data in the optimal feature space.

5 CONCLUSIONS

In this paper, we presented new adaptive algorithms for computation of $\Sigma^{-1/2}$ and introduced a cost function related to them. We proved the convergence of the proposed algorithms using the continuity and initial conditions of the cost function. Simulation results on two class Gaussian data demonstrated the performance of the proposed algorithms for extracting optimal class separability features. The experimental results show that these new adaptive algorithms can be used in many fields of on-line application such as feature extraction for face and gesture recognition. Existence of the cost

function and adaptive nature of the proposed algorithm, make it appropriate to implement related neural networks for different real time application.

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