

ANN-BASED MULTIPLE DIMENSION PREDICTOR FOR SHIP ROUTE PREDICTION

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Abstract: This paper presents a new multiple dimension predictive model based on the diagonal recurrent neural networks (PDRNN) with a combined learning algorithm. This method can be used to predict not only values, but also some points in the multi-dimension space. And also its applications in data mining will be discussed in the paper. Some analysis results show the significant improvement to ship route prediction using the PDRNN model in database of geographic information system (GIS).

1 INTRODUCTION

The problem of prediction is denoted to estimate the output of future according to input and output of now and past in some system. Since Kolmogorov presented a linear optimal predictor in 1941, different kinds of trend analysis methods and prediction models have been used for forecasting and control. In this field, the time series prediction model (Box and Jenkins, 1970) and the self-tuning predictor (Wittenmark, 1974) were two kinds of classical prediction methods. The tradition prediction theories based on time series were developed from linear auto recurrent moving average (ARMA) models. And then these theories were extended to nonlinear process. But, if using the tradition predictive theories, it needs to solve the problems: system modelling, parameter estimating, model modifying and trend forecasting on-line.

In order to solve these problems, some intelligent prediction methods were discussed, in which the forward neural networks with BP algorithm were used more popularly. Prediction based on ANN has made an overwhelming impact on many disciplines. But there are some difficulties in prediction, particularly in the prediction of multi-variable and non-steady dynamic process.

Recent years, scholars had done much research, and made some progresses in this filed. We have researched predictive models using neural networks, such as an ANN-based nonlinear time series model for fault detection and prediction in marine system (Tang, 1998) and an adaptive predictor based on a recurrent neural network for fault prediction and incipient diagnosis (Tang, 2000). Furthermore a direct multi-step adaptive predictor based on a

diagonal recurrent neuron network was presented for intelligent system monitoring (Dou, 2001). These models increased the precision and self-adaptation of prediction in a manner.

However, there existed a problem: former prediction methods based on time series models could only approach or predict processes with one kind of attribute, such as temperature, pressure and flow in an industry process, or stock values and GDP in the economic process. In this case, every parameter must be separately denoted if using a traditional time series model in the dynamic process. But some objects have more than two kinds of attributes, and must be represented as one predictive model. For example, a ship route has two kinds of attributes: longitude and latitude. A satellite position has three kinds of attributes: longitude, latitude and altitude. So the question of how to predict objects with several attributes is an important problem in practice.

This paper discusses self-adaptation prediction methods based on ANN, and presents a multi-dimension predictive model based on parallel diagonal recurrent neuron network (PDRNN) with TD-DBP combined algorithms for time series multi-step forecasting. The paper takes a step forward to use this model in data mining of GIS. Some simulation resolves show the model is able to predict a ship's route according to its position from GPS.

2 PRINCIPLE OF ANN-BASED PREDICTOR

The basic issue of a predictor can be described as: if the past output value series $\{x_t\}$ is known, then

try to design a predictor to obtain the future output value of forward d -step x_{t+d} under the condition of the minimum predictive errors. If x_{t+d} is expressed as $\hat{x}_t(d)$, the model of predictor can be described as follows:

$$\hat{x}_t(d) = f(X_{t-1}, t, P) \quad (1)$$

Where $X_{t-1} = [x_1, x_2, \dots, x_{t-1}]^T$ is the past output value vector, i.e. historical data.

$f(\cdot)$ is a certain nonlinear function.

P is the parameter set of the system model.

The predictor is called the minimum covariance optimal predictor, because the covariance of the predictive error is used for criterion function J as follows:

$$J = \text{Var}(x_{t+k} - \hat{x}_t(k)) \rightarrow \min \quad (2)$$

2.1 RNN-based One-step Predictor

The ANN-based models could be used to construct an adaptive optimal predictor for model identification, parameter correction and value prediction. Assuming that a class of nonlinear autoregressive moving average (NARMA) model, the NARMA (p, q) model is written as:

$$x_t = g(x_{t-1}, x_{t-2}, \dots, x_{t-p}, e_{t-1}, \dots, e_{t-2}, \dots, e_{t-q}) + e_t \quad (3)$$

Where $g(\cdot)$ is an unknown smooth function, and it is assumed that $E(e_t | x_{t-1}, x_{t-2}, \dots) = 0$ and e_t has a finite variance $\text{Var}(e_t) = \sigma^2$. In this case, an approximated condition mean predictor based on the finite past of observations is given by

$$\hat{x}_t = g(x_{t-1}, x_{t-2}, \dots, x_{t-p}, \hat{e}_{t-1}, \hat{e}_{t-2}, \dots, \hat{e}_{t-q}) \quad (4)$$

Where $\hat{e}_j = x_j - \hat{x}_j \quad j = t-1, t-2, \dots, t-q$

For NARMA (p, q) modelling and predicting, a recurrent neural network was presented (Connor, 1994). The recurrent network topology is shown in figure 1. This recurrent neural network (RNN) can be used to approximate the NARMA (p, q) model. The output of the basic RNN-based predictor is

$$\hat{x}_t = \mathbf{W}_o \mathbf{f}(\mathbf{s}) \quad (5)$$

$$\mathbf{s} = \mathbf{W}_{ih} \mathbf{x} + \mathbf{W}_{eh} \mathbf{e} + \boldsymbol{\theta} \quad (6)$$

Where $\mathbf{f}(\cdot)$ is a Sigmoid function vector or other finite continuous monotonically increasing function vectors;

\mathbf{s} is a state vector of the hidden layer;

\mathbf{W}_{ih} is the weight matrix between the input layer and the hidden layer;

\mathbf{W}_{eh} is the weight matrix from feedback units to hidden units;

\mathbf{x} is the input vector;

\mathbf{e} is the error vector;

$\boldsymbol{\theta}$ is a threshold vector.

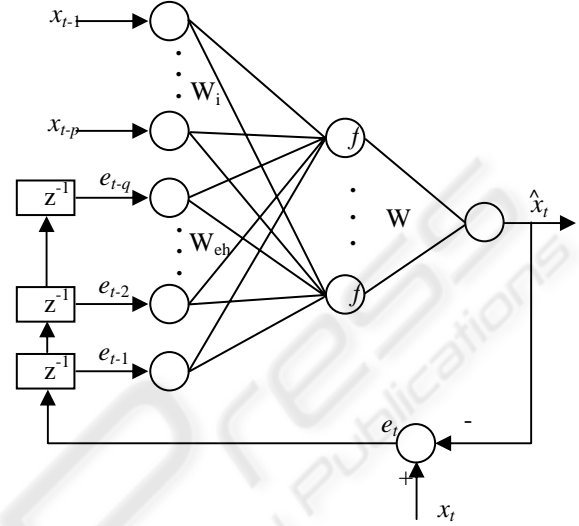


Figure 1: Recurrent network for NARMA models

The parameters of \mathbf{W}_o , \mathbf{W}_{ih} and \mathbf{W}_{eh} are estimated by a dynamic BP learning algorithm (Williams, 1990). That is by learning of RNN to minimize the following error function:

$$E = \frac{1}{2} \sum_{t=p+1}^N (x_t - \hat{x}_t)^2 \quad (7)$$

But, just one step prediction will make by the basic predictor. So some improvement RNN-based predictors had been discussed (Tang, 2000).

2.2 RNN-based Multi-step Predictor

In order to implement the multi-step prediction, the NARMA model should be extended to:

$$\hat{x}_t(d) = g(\hat{X}_t(d-i), \hat{e}_{t+d-j}) \quad (8)$$

Where

$$\hat{e}_{t+d-j} = x_{t+d-j} - \hat{x}_t(d-j), \quad j = 1, 2, \dots, q$$

$$\hat{X}_t(d-i) = [\hat{x}_t(d-1), \hat{x}_t(d-2), \dots, x_t, x_{t-1}, \dots, x_{t-p}]^T$$

The output of the multi-step RNN model is

$$\hat{\mathbf{x}}_t = \mathbf{W}_o \mathbf{f}(\mathbf{s}) \quad (9)$$

Where $\mathbf{s} = \mathbf{W}_{ih} \mathbf{x} + \mathbf{W}_{eh} \mathbf{e} + \boldsymbol{\theta}$

$$\hat{\mathbf{x}}_t = [\hat{x}_t \ \hat{x}_{t+1} \ \cdots \ \hat{x}_{t+d}]^T$$

And the error function is

$$E = \frac{1}{2} \sum_{t=p+1}^N (\mathbf{x}_t - \hat{\mathbf{x}}_t)^T (\mathbf{x}_t - \hat{\mathbf{x}}_t) \quad (10)$$

Using the dynamic BP learning algorithm, and assuming that the dimensions of the input, error, hidden and output matrices could be represented as i , e , h , and o , the iteration formulae of the weight values of the RNN prediction model can be obtained as following:

$$\mathbf{W}(k+1) = \mathbf{W}(k) - \eta \frac{\partial E}{\partial \mathbf{W}} \quad (11)$$

To the weight values of the output layer, there is

$$\frac{\partial E}{\partial \mathbf{W}_o} = - \sum_{t=p+1}^N \mathbf{I}^{(h \times o)} \mathbf{f}(\mathbf{s}) \mathbf{e}_t \quad (12)$$

$$\text{Where } \mathbf{e}_t = \mathbf{x}_t - \hat{\mathbf{x}}_t \quad (13)$$

And \mathbf{I} is an identity matrix.

To the weight values between the hidden layer and input layer, there is

$$\frac{\partial E}{\partial \mathbf{W}_{ih}} = - \sum_{t=p+1}^n (\mathbf{I}^{(i \times h)} \cdot (\mathbf{I}^{(h)} \otimes \mathbf{x}) + \frac{\partial \mathbf{e}}{\partial \mathbf{W}_{ih}} \mathbf{W}_{eh}^T) \cdot \frac{\partial \mathbf{f}(\mathbf{s})}{\partial \mathbf{s}} \cdot \mathbf{W}_o^T \mathbf{e}_t \quad (14)$$

Where \otimes is Kronecker product.

$$\frac{\partial \mathbf{e}}{\partial \mathbf{W}_{ih}} = \left[\frac{\partial e_{t-1}}{\partial \mathbf{W}_{ih}}, \frac{\partial e_{t-2}}{\partial \mathbf{W}_{ih}}, \dots, \frac{\partial e_{t-q}}{\partial \mathbf{W}_{ih}} \right] \quad (15)$$

To the weight values of the hidden layer and feedback units, there is

$$\frac{\partial E_t}{\partial \mathbf{W}_{eh}} = - \sum_{t=p+1}^N \left(\mathbf{I}^{(e \times h)} (\mathbf{I}^{(h)} \otimes \mathbf{e}) + \frac{\partial \mathbf{e}}{\partial \mathbf{W}_{eh}} \mathbf{W}_{eh}^T \right) \frac{\partial \mathbf{f}(\mathbf{s})}{\partial \mathbf{s}} \mathbf{W}_o^T \mathbf{e}_t \quad (16)$$

Where

$$\frac{\partial \mathbf{e}}{\partial \mathbf{W}_{eh}} = \left[\frac{\partial e_{t-1}}{\partial \mathbf{W}_{eh}}, \frac{\partial e_{t-2}}{\partial \mathbf{W}_{eh}}, \dots, \frac{\partial e_{t-q}}{\partial \mathbf{W}_{eh}} \right] \quad (17)$$

3 DRNN PREDICTIVE MODELS

In order to obtain the optimal predictive value of the future output of the analyzed system based on its historical data, a stochastic dynamic model of the

analyzed system should be set up, which can modify the model parameter adaptively. A diagonal recurrent neural network was used to represent the dynamic process based on NARMA model (Dou, 2001). The NARMA model is defined as:

$$\hat{x}_{t+d} = h(x_{t-1}, x_{t-2}, \dots, x_1, \hat{e}_{t-1}, \dots, \hat{e}_1) \quad (18)$$

Where $\hat{e}_j = x_j - \hat{x}_j$, $j = t-1, t-2, \dots, 1$;

$h(\cdot)$ is a nonlinear function.

Figure 2 shows the structure of the DRNN-based predictive model. The neural network model with two inputs and several outputs includes three layers. In order to realize direct multi-steps prediction, the output layer composes of d linear neural units. And the middle layer (i.e. hidden layer) makes up NH nonlinear dynamic neurons whose map function is the sigmoid function, and each of the hidden unit includes a self-feedback with one step delay (recursion layer). The input layer includes two linear neurons, and one of them accepts x_{t-1} as input signal, another accepts \hat{x}_{t-1} , which is one-step delay of the output x_t . This network can be regarded as a parsimonious version of the Elman-type network. It has a diagonal structure, that is, there is no interaction between different dynamic neurons.

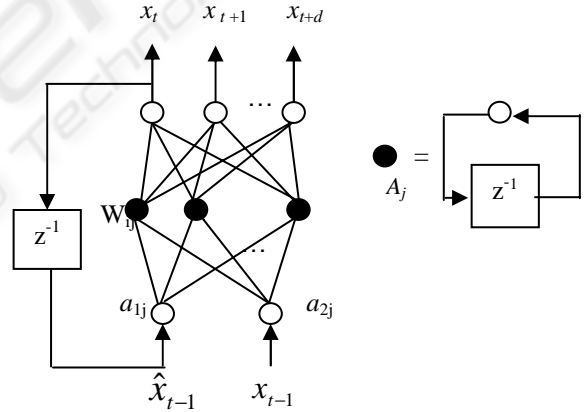


Figure 2: DRNN-based predictive model

The transfer function of this network is described as follows. Suppose w_{ij} ($i=1, 2, \dots, NH$; $j=1, 2, \dots, d$) are the connection weights of the output layer, A_i ($i=1, 2, \dots, NH$) are the connection weights of the recursion layer, a_{1i} ($i=1, 2, \dots, NH$) are the connection weights of the input x_{t-1} to each of the hidden units, a_{2i} ($i=1, 2, \dots, NH$) are the connection weights of the input \hat{x}_{t-1} to each of the hidden units, the output of the neural unit of the output layer is expressed as $\hat{x}_t(j)$, ($j=1, 2, \dots, d$), the output of an neurons of hidden layer is expressed as $H_i(t)$, ($i=1, 2, \dots, NH$), the input of the neural unit of the hidden layer is expressed as $V_i(t)$ ($i=1, 2, \dots$,

NH), then the map relations of the DRNN are shown as follows:

$$\hat{x}_t(j) = \sum_{i=1}^{NH} [w_{ij}H_i(t) - \theta_{oj}] \quad (19)$$

$$\text{Where } H_i(t) = g_N(V_i(t) - \theta_{hi}) \quad (20)$$

$$V_i(t) = A_i H_i(t-1) + a_{hi}(x_{t-1} - \theta_1) + a_{2i}(\hat{x}_{t-1} - \theta_2) \quad (21)$$

$g_N(\cdot)$ is sigmoid function, θ_{oj} is the threshold of the neural unit of output layer, θ_{hi} is the threshold of the neuron of hidden layer. The initial conditions of this model are $H_i(0) = 0$ and $V_i(0) = 0$. So the transfer function of the neural network is a nonlinear continuous function, and the output of the neural network $\hat{x}_t(j)$ is an appropriate nonlinear function of all input signals (x_1, x_2, \dots, x_{t-1}).

It is known that the supervised learning algorithm based on the error between the actual output and the anticipant output is not suitable for the direct multi-step predictive model. And a neural network with the fixed structure and parameters is difficult or even impossible to express the inherent dynamic performance of the uncertain nonlinear systems. For this reason the temporal difference (TD) learning algorithm and the dynamic back-propagation (DBP) algorithm are synthesized for the network training. The combined learning algorithm will adaptively modify the parameters of the predictive model according to the errors between the predictive value and the actual detective value.

Suppose P_j^t is the output of the j th output neuron at t time, i.e. $\hat{x}_t(j)$, P_{j-1}^{t+1} is the output of the $(j-1)$ th output neuron at $t+1$ time, i.e. $\hat{x}_t(j)$. It is obvious that P_j^t and P_{j-1}^{t+1} are the predictive output value of the analyzed system at the same time. And they should be equal if the prediction is accurate. So the P_{j-1}^{t+1} can be used as the expectant output of the P_j^t . This is the basis of TD learning algorithm. The training error of one learning sample of the DRNN is expressed as:

$$e = \frac{1}{2} \sum_{j=0}^K (x_{t+j} - \hat{x}_t(j))^2 \quad (22)$$

According to TD learning algorithm, it can be expressed as

$$e = \frac{1}{2} \sum_{j=0}^K (P_j^{t+1} - \hat{x}_t(j))^2 \quad (23)$$

Here P_0^{t+1} is defined as the expectant output of x_t at t time.

4 A MULTI-DIMENSION PREDICTOR BASED ON PDRNN

One dimension predictive model can predict an object with one kind of attribute. A multi-dimension predictive model can predict an object with more than one kind of attribute at the same time. In the multi-dimension predictive model, there are different relations in different attributes and the relations can be changed by a dynamic process. For this reason, ANN-based adaptive predictors must be introduced to modify the parameters of a predictive model on-line.

4.1 The Framework of PDRNN Model

In order to solve the predictive problem of objects with multi-attributes, this paper presents a new multi-dimensional predictive model based on the diagonal recurrent neural networks (PDRNN) with a parallel combined learning algorithm. Fig. 3 shows the framework of PDRNN model. There are four layers in this model, the first layer is the network input layer; the second layer is the network input assignment layer; the third layer is the network hidden layer, in which every hidden unit includes a self-feedback with one step delayed; the forth layer is the network output layer, (the network output layer connects with the network input layer through one-step feedback). There are n dimension variables to input paralleled in the network input layer. It solves the problem that only one variable can be input in some one dimension predictive models, and the PDRNN model can predicts an object with multiple variables and attributes.

4.2 The Mathematical Description of the PDRNN Model

As shown in figure 3, there are p input units in the network input layer, every input unit has n dimension variables, X_t can be obtained by Z_{t+1} with one step delayed, so each attribute variable has $p-1$ input values in this network every time in fact. The network input assignment layer assigns n values of each variable to n sub-input layers paralleled, as shown in equation (24). The hidden layer has n sub-hidden layers paralleled, every sub-hidden layer has

NH^1 (the number of sub hidden layer) nonlinearity units with S functions or T functions, the value of NH^1 can be changed, and every sub-hidden layer has self-feedback with one step delayed. In this network, n paralleled sub-networks consisted of the sub-input layers and sub-hidden layers, all the parallel sub-networks respectively train different attributes at the same time. This network neglected the relations in the different attributes and attribute values. There are k linear units in the network output layer, which can do d -step prediction at most. The mathematical model could be described as follows:

The network input layer

$$[X_t, X_{t-1}, \dots, X_{t-p}]$$

Where $X_t = [X_t^1, X_t^2, \dots, X_t^n]$ (24)

The I th parallel sub-network's sub input layers:

$$[X_t^I, X_{t-1}^I, \dots, X_{t-p}^I] \quad (25)$$

The I th parallel sub-network's sub hidden layers:

$$H^I_i(t) = g_N(V^I_i(t)) \quad (26)$$

In equation (26), at every t time, $H^I_i(t)$ is the output of the i th hidden unit in the I th parallel sub-network, $g_N(\cdot)$ is Sigmoid (S) function or Tangent (T) function. And there:

$$V^I_i(t) = A^I_i H^I_i(t-1) + \sum_{i=1}^{N^I} w^I_i X^I_{t-1} \quad (27)$$

In equation (27), at every t time, $\sum_{i=1}^{N^I} w^I_i X^I_{t-1}$ is the sum of all the inputs in the i th hidden layer of the I th parallel sub-network, w^I_i is the normalized (relative) fulfilment weights between the sub-input layer and sub-hidden layer of the I th parallel sub-network, A^I_i is the self-recurrent layer's relative weights of the i th hidden unit in the I th parallel sub-network. Z_{t+j} is defined as the output of the j th output unit and includes all the attribute values at this time. So the output variables of the network model could be written by equation (28) as below:

$$Z_{t+j} = \left(\sum_{i=1}^{NH^1} a^1_{ij} H^1_i(t), \sum_{i=1}^{NH^2} a^2_{ij} H^2_i(t), \dots, \sum_{i=1}^{NH^n} a^n_{ij} H^n_i(t) \right) \quad (28)$$

Where $Z^I_{t+j} = \sum_{i=1}^{NH^I} a^I_{ij} H^I_i(t)$

is the j th output layer's output of the I th attribute at t time. In the initialization, the threshold value of all the nerve units were neglected at every t time, and

$$H^I_i(0) = 0$$

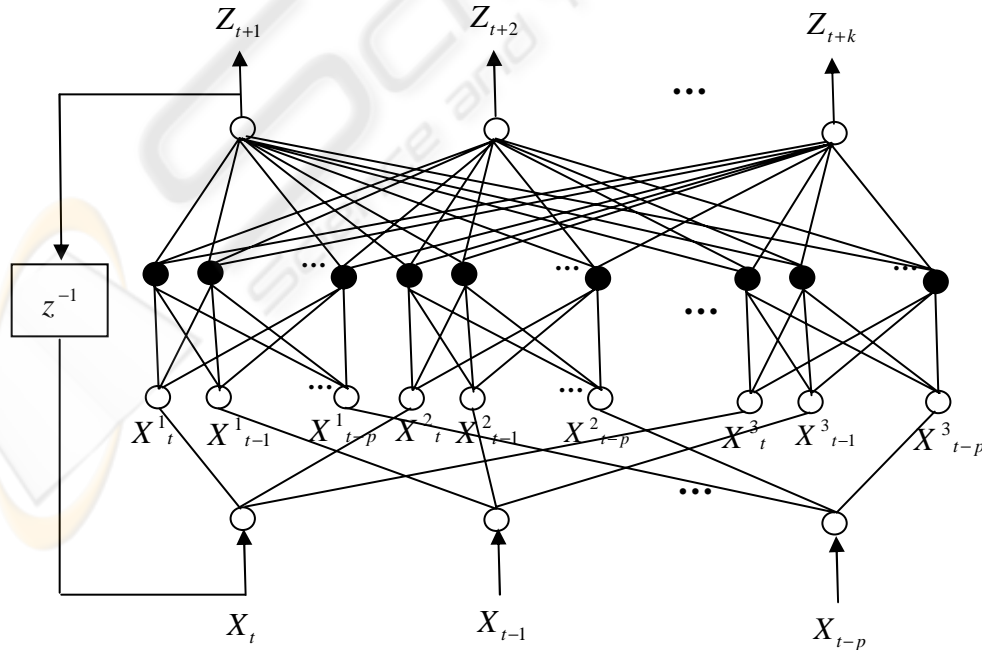


Figure 3: Multi-dimension predictive model based on PDRNN

5 THE LEARNING ALGORITHM

This paper combines the time different method (TD) and dynamic BP method to train PDRNN model. If Z^I_{t+j} as the j th output of the I th attribute at t time, is the predictive value of the I th attribute at $t+j$ time, X^I_{t+j} is real value of the I th attribute at $t+j$ time in the future.

In the ideal condition, Z^I_{t+j} is equal to X^I_{t+j} , but usually there are some errors in practice. It could be represented by following error function.

$$e^I = \frac{1}{2} \sum_{j=0}^{N^I_k} (X^I_{t+j} - Z^I_{t+j})^2 \quad (29)$$

Equation (29) is the training error of the I th attribute at $t+j$ time. Here use a dynamic BP method to correct relative weighs. The learning algorithm is as follows

$$\frac{\partial e^I}{\partial a^I_{ij}} = \frac{\partial e^I}{\partial X^I_{t+i}} \frac{\partial X^I_{t+i}}{\partial a^I_{ij}} = -(Z^I_{t+j} - X^I_{t+j}) H^I_i(t) \quad (30)$$

Where a^I_{ij} are the normalized (relative) fulfilment weights between the sub hidden layer and sub output layer of the I th paralleled sub-network.

The formula corrected of a^I_{ij} as follows:

$$\begin{aligned} a^I_{ij}(t+1) &= a^I_{ij}(t) - \xi^I \frac{\partial e^I}{\partial a^I_{ij}} \\ &= a^I_{ij}(t) + \xi^I (X^I_{t+j} - Z^I_{t+j}) H^I_i(t) \end{aligned} \quad (31)$$

Where ξ^I is learning parameter of the I th attribute's output layer, w^I_{ij} is the normalized (relative) fulfilment weights between the sub input layer and sub hidden layer of the I th paralleled sub-network.

$$\begin{aligned} \frac{\partial e^I}{\partial w^I_{ij}} &= - \sum_{j=0}^{N^I_k} \frac{\partial e^I}{\partial X^I_{t+i}} \frac{\partial X^I_{t+i}}{\partial w^I_{ij}} \frac{\partial H^I_i(t)}{\partial w^I_{ij}} \\ &= - \sum_{j=0}^{N^I_k} (X^I_{t+j} - Z^I_{t+j}) w^I_{ij} \frac{\partial H^I_i(t)}{\partial w^I_{ij}} \end{aligned} \quad (32)$$

The formula is corrected for w^I_{ij} as follows:

$$\begin{aligned} w^I_{ij}(t+1) &= w^I_{ij}(t) - \eta^I \frac{\partial e^I}{\partial w^I_{ij}} \\ &= w^I_{ij}(t) + \eta^I \sum_{j=0}^{N^I_k} (X^I_{t+j} - Z^I_{t+j}) a^I_{ij} Q^I_{ij}(t) \end{aligned} \quad (33)$$

Where η^I is learning parameter of the I th attribute's input layer;

A^I_i are the self-recurrent layer's relative weights.

$$\begin{aligned} \frac{\partial e^I}{\partial A^I_i} &= - \sum_{j=0}^{N^I_k} \frac{\partial e^I}{\partial X^I_{t+i}} \frac{\partial X^I_{t+i}}{\partial H^I_i(t)} \frac{\partial H^I_i(t)}{\partial A^I_i} \\ &= - \sum_{j=0}^{N^I_k} (X^I_{t+j} - Z^I_{t+j}) a^I_{ij} P^I_i(t) \end{aligned} \quad (34)$$

Where $P^I_i = \frac{\partial e^I}{\partial A^I_i}$, and $P^I_i(0) = 0$.

The formula corrected of A^I_i as follows:

$$\begin{aligned} A^I_i(t+1) &= A^I_i(t) - \mu^I \frac{\partial e^I}{\partial A^I_i} \\ &= A^I_i(t) + \mu^I \sum_{j=0}^{N^I_k} (X^I_{t+j} - Z^I_{t+j}) a^I_{ij} P^I_i(t) \end{aligned} \quad (35)$$

Where μ^I is the learning parameter of the I th attribute's self-recurrent layer.

First is to adjust the network framework, make sure of the number of input layers, hidden layers and output layers, and the number of maximal learning steps. The above parameters are significant for predictive accuracy. Second initialize the network (for new data, the initialization is random), then ANN begins learning. Equation (29) serves as a standard to judge the predictive value of each attribute. In order to avoid the learning of ANN falling into a dead area, the learning step could be adjusted according to different attribute values. Here a " \bar{e} " function was presented, equation (36) serves as a standard to judge the predictive values of all the attributes.

$$\bar{e} = \frac{1}{n-k} \sum_{t=k}^n E_t \quad (36)$$

$$\text{And } E_t = \sqrt{\sum_{l=1}^m (e^I_t)^2} \quad (37)$$

Where e^I_t is the error between the real value and the predictive value of the I th attribute at t time, E_t can be changed according to different attributes. \bar{e} is the average error, used to judge the convergence of the multi-dimension predictive model based on PDRNN. Because the initialization of ANN is random, the first couple of predictive values may be not very good, \bar{e} is computed from the k th predictive value. The training will be stopped when

\bar{e} is up to standard, or reset up the network framework until \bar{e} confirms to requirements.

6 SIMULATIONS AND APPLICATION

The prediction based on PDRNN extends ANN-based time series prediction model from a single attribute to a multi-attribute. Figure 4 is a three-step predictive value contrasting with real values of straight line, real lines are real values, “*” indicates the predictive values. The points of straight line have two kinds of attributes, every parallel sub-network has ten sub input layers and fifteen sub hidden layers. T function is used in the hidden units, the maximum of e is $2.265e^{-5}$, \bar{e} is $8.36e^{-8}$.

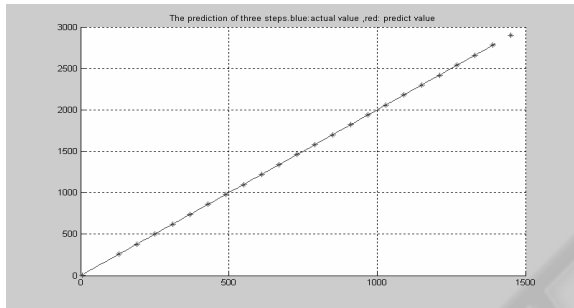


Figure 4: Predictive value contrasting

Figure 5 is a three-step predictive value of nonlinearity as follows:

$$y = \begin{cases} 2 \times x, & 0 \leq x < 2\pi \\ \sin(x), & 2\pi \leq x < 4\pi \end{cases}$$

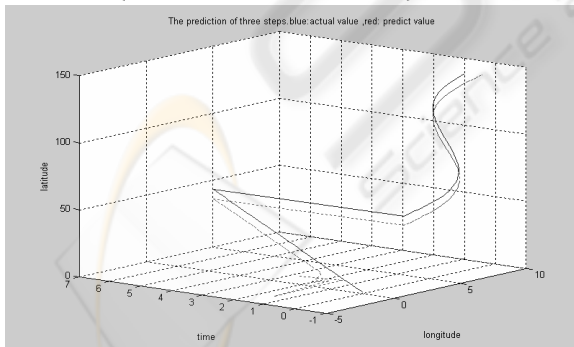


Figure 5: Prediction of nonlinearity

Figure 6 is a three-step predictive value of 3D curve contrasting with real value, T function is used in the hidden units, every parallel sub-network only has ten sub input layers and fifteen sub hidden layers. The learning step of 3D curve is more than 2D curve.

The maximum predictive error of 3D curve is 0.0011, \bar{e} is $8.265e^{-4}$.

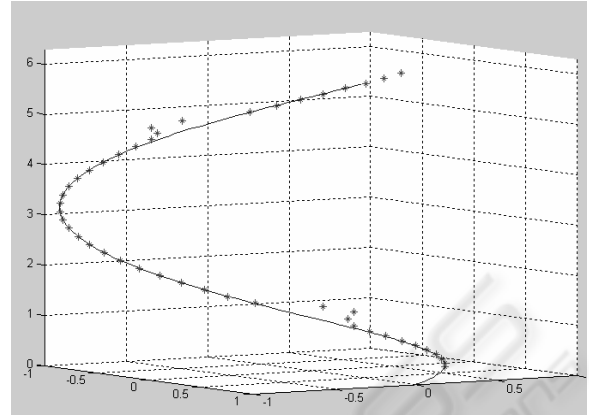


Figure 6: Prediction of 3D curve

An application process for GIS in Marine Engineering with the predictor based on PDRNN is shown in figure 7.

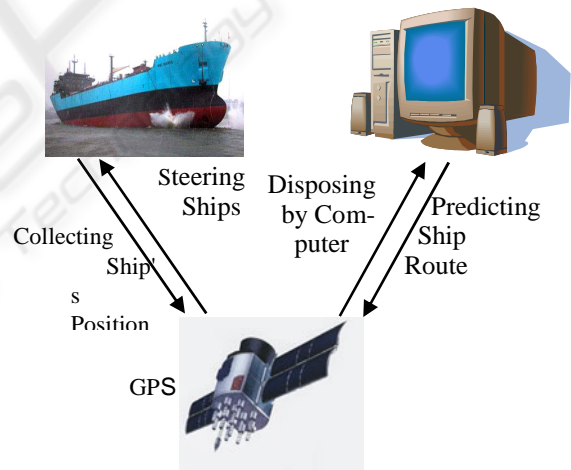


Figure 7: GIS in Marine Engineering

In the system, data recorded abundant ship positions from GPS, and established a database through ACCESS. The software of GIS is ArcView 3.2 of ESRI (Goodchild, 1992). ArcView 3.2 called the data from ACCESS, and eliminated unnecessary data by applying ArcView 3.2. After data pre-processing, the ship route could be selected and drawn in an electronic chart. Using the PDRNN predictor the tracking of the ship route could be forecasted.

The process of ship route prediction by means of a PDRNN model is as follows: first, get the distribution data of the ship's position from GPS (Wang, 2003), then select sample points. In this

paper, the each sample point was selected every 2.5 hours. Figure 8 is a selected ship route. Finally, the ship route was predicted by means of a PDRNN model.

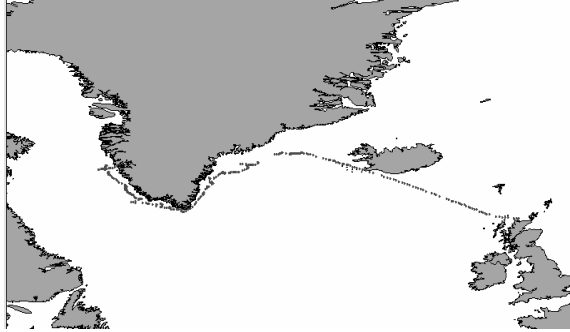


Figure 8: The ship route in GIS

Figure 9 is another prediction resolve of a ship route. This chart shows that this ship route is a variant random process, but the predictive algorithm based on a PDRNN model can follow this process, and do three-step prediction. The predictive maximum error of the ship route is $1.065e^{-14}$, \bar{e} is $3.326e^{-15}$. Thus the error of prediction is small.

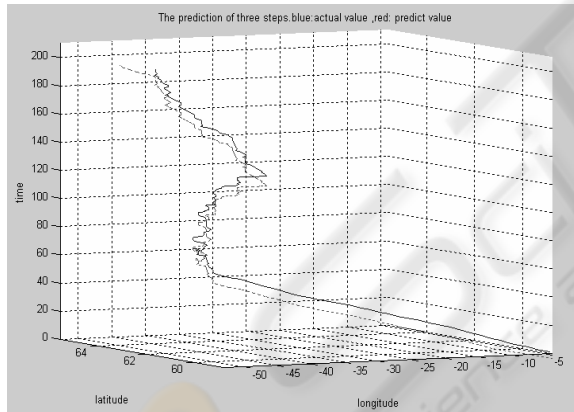


Figure 9: The prediction of ship route

7 CONCLUSIONS

As mentioned above, the NARMA models based on the recurrent neural networks with a dynamic BP algorithm is suited for trend prediction. This paper presents a multi-dimension predictive model based on PDRNN. This predictor could store in memory all the past input information, network output is some nonlinear function of all the past input, and this model can realize a nonlinear dynamic mapping. As a predictive model, the framework is very simple, the dynamic behavior of this model could

regulate network frameworks with self-adaptation, and this model could predict an object with multi-attribute. Moreover, this paper has presented an “ \bar{e} ” function, which can judge if the whole network structure confirms to requirements, and has presented an “input plus” method, which can reduce the training time. The application in ship routing prediction shows the new predictive model is better to predict a multiple dimension dynamic process.

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