Prediction of Reagents Needs Using Radial Basis Function in Teaching Hospital

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Abstract—A robust reagents prediction is able to support the service improvement in laboratories. In this paper, Radial Basis Function Networks (RBFN) method with (3, Q, 1) architecture is used to predict two types of reagents needs, i.e. SD Bioline HBsAg and SD Bioline Anti HCV. Data of reagents from 2012 - 2013 are used as training data, whereas 2014 data are used as comparative data for the prediction result. In RBFN training, the best condition obtained when the spread value is 4 with RMSE 1.63E-06 for both types of reagents. The prediction results with RBFN methods reached 99% with correlation value of 0.99 for each reagents. RBFN method shows better prediction result compared to BPNN method with prediction of 92%.

Keyword-Reagents prediction, Radial basis function network, Teaching hospital, BPNN, RMSE

I. INTRODUCTION

A laboratory installation is one operational support unit in hospital with the duty and responsibility in patient's illness examination as well as one of the service unit that can be developed into a costs center and revenues for the hospital. The problems are often occurred in a laboratory installation is a use of budget for the reagents purchasing. It is happen due to the lack of knowledge of the laboratory personnel to analyzing budget needs so the number of reservations is mismatched with the actual reagents needed at the laboratory. By using empirical data at the hospital, these prediction errors can actually be overcome by using appropriate artificial intelligence methods with the data condition to be predicted.

In several existing research, wide variety of theories and methods for the needs logistics prediction in hospital or the prediction of reagents needs have been carried out. Cao et. al. studied the hospital logistics stock demands forecasting based on data mining and back propagation neural network. The result accuracy of this research is 87% [1]. In addition, Wen et al. examined the model of hospital engineering logistics management with JIT theory aimed at ensuring the logistics operations can run normally and reduce excessive operational costs [2]. While research on the prediction of reagents needs have been described as Decision Support System mas main tools [3-5]. Anon et al. doing research on decision-making model development for reagents forecasting with Analysis Hierarchy Process (AHP) method, Goal Programming (GP), Time Series Analysis (TSA) and Expert Systems (ES). The results of this study showed that the ES method provide better results than other methods with an average MSE of 3.75 [3]. Furthermore, Tzu et al. examines the decision-making management for reagents purchase with exponential smoothing method. This study obtained different results for each type of reagents with each correlation values for reagents plateau type is 0.3-0.5 and the increment type and decrement reagents type correlation value is 0.5-0.7 [4].

Baizul et.al also made safety stock reagents prediction with Back Propagation Neural Network (BPNN), where the accuracy level of the prediction results reached 92% [5]. It is proven that shifting error learning from back propagation to radial basis function will increase accuracy of system not only on data series but also for multimedia matrices [6,7]. Based on the prediction results in previous studies, the accuracy level of predicted results still need to be improved. Therefore, Radial Basis Function Network (RBFN) method is used in this study to optimize prediction accuracy rate.

II. LITERATURE REVIEW

A. Radial Basis Function Network

Radial Basis Function Network (RBFN) is an artificial neural network model which transform input nonlinearly using radial basis activation function similar to the Gaussian activation function in hidden layer unit before linearly processed on the output layer. RBFN consists of three layers, i.e. input layer, hidden layer and output layer. Each hidden layer represents activation function in the form of radial basis function. Each input of its network will enable all the activation function on hidden layer. Each unit on hidden layer is called basis functions. Hidden layer consists of a number of similar basis functions that will generate an output with certain weight. The output of this network is the sum of all basis functions output multiplied by its weight [6].

B. Proposed RBFN Architecture

In this paper, RBFN architecture have three input layer, hidden layer contain neurons as much as the data input and one output layer that uses only one neuron (3, Q, 1) that is shown in Fig. 1. Each input activates each function base on its own network. For example, input x_1 will activate basis function on the first RBFN network, subsequently basis functions φ_{11} , φ_{12} and φ_{13} be activated. Furthermore, cross-correlation between each basis functions on each network is calculated and the results are weighted w_{ij} . RBFN network output is calculated by the following equation [6].

$$Output = W_0 + \sum_{i=1}^{3} \sum_{j=1}^{3} \varphi_{ij} W_{ij}$$
(1)

Where w_0 is bias weight, φ_{ij} is the activation basis function which is defined by Gaussian function as follows.

$$\boldsymbol{\varphi}_{ij} = \boldsymbol{e}^{\left(-\frac{\|\boldsymbol{x}-\boldsymbol{c}_{ij}\|^2}{\sigma_{ij}^2}\right)}$$
(2)

with c_i is the jth gaussian functions center, σ_i is the jth gaussian function width and x is the basis function input.

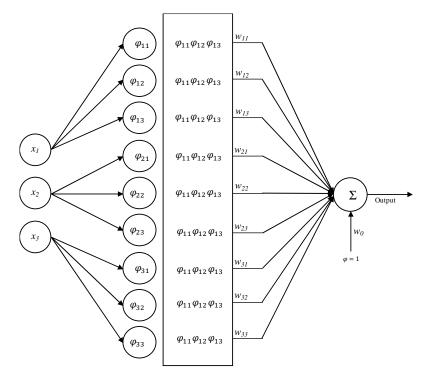


Fig 1. Proposed RBFN Architectural Design for Reagents Input

III. METHODOLOGY

A. Preprocessing

Data reagents in this paper are collected from the laboratory unit of the teaching hospital for 3 years, from 2012 to 2014. These data consists of three variables, i.e. reagents consumption, reagents demand and residual reagents. The data from 2012 to 2013 is used as training data while the data in 2014 is used validation data. At the initial stage, the data is normalized before the training process by using the following equation.

$$x' = (x-b)/a - b$$

where x' is the normalized data, x is the input data, a represent the largest data value and b is the smallest data value.

B. Training Process

Training process consists of five steps, i.e entering the input variables, normalizing the data, initializing network, setting the spread value and calculating the error value. Spread value setting is performed to optimize network activation function. The training process will stop if the error value < 0.00001. The network training process of RBFN is shown in Fig. 2.

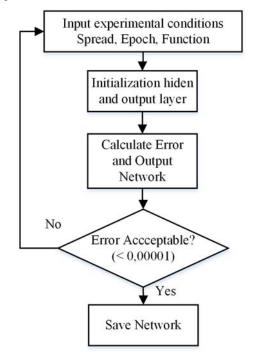


Fig 2. The network training of RBFN

C. Postprocessing

At this stage, the training output will be denormalized with the following equation.

$$p' = y(a-b) + b$$

(4)

(3)

where p' is the denormalized data, y is the output of the training, b is the smallest data and a is the largest data. D. Process Validation and Prediction

Validation process are carried out in several stages, such as inserting test data of year 2014 that consists of three variable parameters (residual reagents data, reagents consumption data, reagents demand data), data normalization process to obtain data interval from 0 to 1, input the best network (*net*) from RBFN training result, conducting process of reagents needs prediction for 2014, perform data denormalization to restore the data to the actual value and calculate correlation between the prediction results and the actual laboratory reagents.

Several steps being taken in the validation process are entering test data in 2014, normalizing data, input the best net value of the training results, conducting process of reagents needs prediction for 2014, denormalized data and calculate the correlation of the prediction results and the actual data.

IV. RESULTS AND DISCUSSION

A. Result Training Network

In this paper, the best training results are based on the smallest RMSE values that influenced by the spread value. The network training results are shown in Table I.

	Spread Value	RMSE Value		
No		SD Bioline Reagents HbsAg	SD Bioline Reagents Anti HCV	
1	2	3.25E-06	2.31E-06	
2	4	1.63E-06	1.63E-06	
3	6	9.62E-06	2.82E-06	
4	8	6.71E-06	4.31E-06	
5	10	3.64E-05	3.98E-06	
6	12	2.60E-05	1.29E-05	
7	14	1.05E-04	4 1.08E-05	
8	16	7.87E-05	1.21E-05	
9	18	6.17E-05	1.28E-04	
10	20	5.25E-05	1.04E-04	
11	22	1.83E-04	1.02E-04	
12	24	1.55E-04	8.78E-05	
13	26	1.35E-04	7.50E-05	
14	28	1.14E-04	6.20E-05	
15	30	1.01E-04	5.49E-05	
16	32	8.96E-05	5.27E-05	
17	34	8.01E-05	6.73E-05	
18	36	7.60E-05 6.06E-05		
19	38	7.45E-05	1.86E-04	
20	40	8.13E-05	1.63E-04	

TABLE I. Network Training of RBFN with adjustment spread value

The best RBFN network training results achieved when the spread value is 4 with RMSE value of 1.63E-06 for both reagents. The RMSE value on training calculated by the following equation [7].

$$RMSE = \frac{\sqrt{\frac{1}{N} \sum_{t=h}^{N} (y_t - y'_t)^2}}{y_{\text{max}} - y_{\text{min}}}$$
(5)

where y_t is the actual *data*, N is the data amount, y_t is the prediction data, y_{max} is the maximum value of the actual data and y_{min} is minimum value of the actual data.

B. Prediction Results Comparison Between RBFN and BPNN Methods

The predicted results accuracy of RBFN method in this paper in the form of a correlation value compared with the predicted results of BPNN is discussed in Ref. [1] as shown in Table II.

Mt	SD Bioline HBsAg			SD Bioline Anti HCV		
	Actual	RBFN	BPNN	Actual	RBF N	BPNN
1	260	260	267	180	180	157
2	670	670	561	100	100	61
3	200	200	182	180	180	193
4	260	260	239	260	260	282
5	200	200	190	100	100	111
6	300	300	273	310	310	348
7	350	350	352	250	250	256
8	430	430	496	380	380	365
9	450	450	369	350	350	361
10	150	150	151	300	300	343
11	300	300	276	250	250	347
12	600	600	550	500	500	367
Crl		0.99	0.96		0.99	0.88

TABLE II. Comparison of Results Prediction Method RBFN with BPNN

Act=Actual, Mt=Target of Month, Crl=Correlation

In Table II, both reagents show the same correlations values of 0.99 for RBFN prediction method. Based on the correlation values in Table 2, it can be said that the results of prediction methods RBFN better than BPNN method. To calculate the correlation value prediction results above used the following equation [8]:

$$r = \frac{n\sum xy - (\sum x)(\sum y)}{\sqrt{\{n\sum x^2 - (\sum x)^2\}\{n\sum y^2 - (\sum y)^2\}}}$$
(6)

Where *n* is the number of data couple *x* and *y*, Σx is the total amount of *x*, Σy is the total amount of *y*, Σx^2 is the square from the total number of *x*, Σy^2 is the square of the total number of *y* and Σxy is the result of the total number of variables *x* and *y* variables multiplication.

As for predictions result comparison graph between RBFN with BPNN method is shown in Fig. 3 and Fig. 4. The results show that the RBFN prediction is more accurate than that of BPNN.

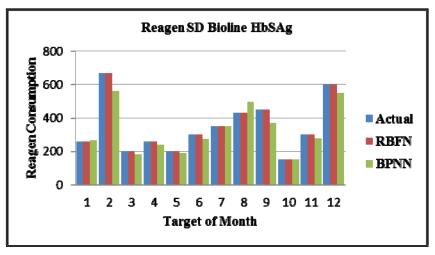


Fig 3. Comparison of Results Prediction Methods RBFN and BPNN For SD Bioline HBsAg Reagents

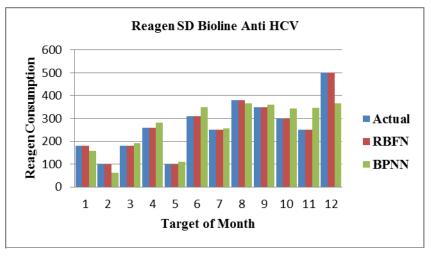


Fig 4. Comparison of Results Prediction Methods RBFN and BPNN For SD Bioline Anti HCV Reagents

Fig. 3 and Fig. 4 shows the results predicted by the RBFN method is quite accurate and shadowing the reagents consumption pattern appropriate with the actual data. The red graph shows the RBFN method prediction, green is BPNN methods prediction and blue is the actual reagents usage data.

V. CONCLUSION

Prediction of reagents SD Bioline HBsAg and SD Bioline Anti-HCV using Radial Basis Function Network (RBFN) method has been studied in this paper. With RBF Architecture method (3, Q, 1) and spread value of 4, the prediction results profound the the correlation of 0.99 and 99% accuracy rate for both types of reagents. Based on the results prediction accuracy, RBFN method shows better results than BPNN method with the prediction results of 92%.

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