



ELSEVIER

European Journal of Operational Research 116 (1999) 16–32

EUROPEAN
JOURNAL
OF OPERATIONAL
RESEARCH

Theory and Methodology

Artificial neural networks in bankruptcy prediction: General framework and cross-validation analysis

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Received 10 March 1997; accepted 22 December 1997

Abstract

In this paper, we present a general framework for understanding the role of artificial neural networks (ANNs) in bankruptcy prediction. We give a comprehensive review of neural network applications in this area and illustrate the link between neural networks and traditional Bayesian classification theory. The method of cross-validation is used to examine the between-sample variation of neural networks for bankruptcy prediction. Based on a matched sample of 220 firms, our findings indicate that neural networks are significantly better than logistic regression models in prediction as well as classification rate estimation. In addition, neural networks are robust to sampling variations in overall classification performance. © 1999 Elsevier Science B.V. All rights reserved.

Keywords: Artificial intelligence; Neural networks; Bankruptcy prediction; Classification

1. Introduction

Prediction of bankruptcy has long been an important topic and has been studied extensively in the accounting and finance literature [2,3,6,16,29,30]. Since the criterion variable is categorical, bankrupt or nonbankrupt, the problem is one of classification. Thus, discriminant analysis, logit and probit models have been typically used for this purpose. However, the validity and effectiveness of these conventional statistical methods

depend largely on some restrictive assumptions such as the linearity, normality, independence among predictor variables and a pre-existing functional form relating the criterion variable and predictor variables. These traditional methods work best only when all or most statistical assumptions are apt. Recent studies in artificial neural networks (ANNs) show that ANNs are powerful tools for pattern recognition and pattern classification due to their nonlinear nonparametric adaptive-learning properties. ANN models have already been used successfully for many financial problems including bankruptcy prediction [62,67].

Many researchers in bankruptcy forecasting including Lacher et al. [33], Sharda and Wilson

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[57], Tam and Kiang [61], and Wilson and Sharda [66] report that neural networks produce significantly better prediction accuracy than classical statistical techniques. However, why neural networks give superior classification is not clearly explained in the literature. Particularly, the relationship between neural networks and traditional classification theory is not fully recognized [51]. In this paper, we provide explanation that neural network outputs are estimates of Bayesian posterior probabilities which play a very important role in the traditional statistical classification and pattern recognition problems.

In using neural networks, the entire available data set is usually randomly divided into a training (in-sample) set and a test (out-of-sample) set. The training set is used for neural network model building and the test set is used to evaluate the predictive capability of the model. While this practice is adopted in many studies, the random division of a sample into training and test sets may introduce bias in model selection and evaluation in that the characteristics of the test may be very different from those of the training. The estimated classification rate can be very different from the true classification rate particularly when small-size samples are involved. For this reason, it is one of the major purposes of this paper to use a cross-validation scheme to accurately describe predictive performance of neural networks. Cross-validation is a resampling technique which uses multiple random training and test subsamples. The advantage of cross-validation is that all observations or patterns in the available sample are used for testing and most of them are also used for training the model. The cross-validation analysis will yield valuable insights on the reliability of the neural networks with respect to sampling variation.

The remainder of the paper will be organized as follows. In Section 2, we give a brief description of neural networks and a general discussion of the Bayesian classification theory. The link between neural networks and the traditional classification theory is also presented. Following that is a survey of the literature in predicting bankruptcy using neural networks. The methodology section contains the variable description, the data used and the design of this study. We then discuss the cross-

validation results which will be followed by the final section containing concluding remarks.

2. Neural networks for pattern classification

2.1. Neural networks

ANNs are flexible, nonparametric modeling tools. They can perform any complex function mapping with arbitrarily desired accuracy [14,23–25]. An ANN is typically composed of several layers of many computing elements called nodes. Each node receives an input signal from other nodes or external inputs and then after processing the signals locally through a transfer function, it outputs a transformed signal to other nodes or final result. ANNs are characterized by the network architecture, that is, the number of layers, the number of nodes in each layer and how the nodes are connected. In a popular form of ANN called the multi-layer perceptron (MLP), all nodes and layers are arranged in a feedforward manner. The first or the lowest layer is called the input layer where external information is received. The last or the highest layer is called the output layer where the network produces the model solution. In between, there are one or more hidden layers which are critical for ANNs to identify the complex patterns in the data. All nodes in adjacent layers are connected by acyclic arcs from a lower layer to a higher layer. A multi-layer perceptron with one hidden layer and one output node is shown in Fig. 1. This three-layer MLP is a commonly used ANN structure for two-group classification problems like the bankruptcy prediction. We will focus on this particular type of neural networks throughout the paper.

Like in any statistical model, the parameters (arc weights) of a neural network model need to be estimated before the network can be used for prediction purposes. The process of determining these weights is called training. The training phase is a critical part in the use of neural networks. For classification problems, the network training is a supervised one in that the desired or target response of the network for each input pattern is always known a priori.

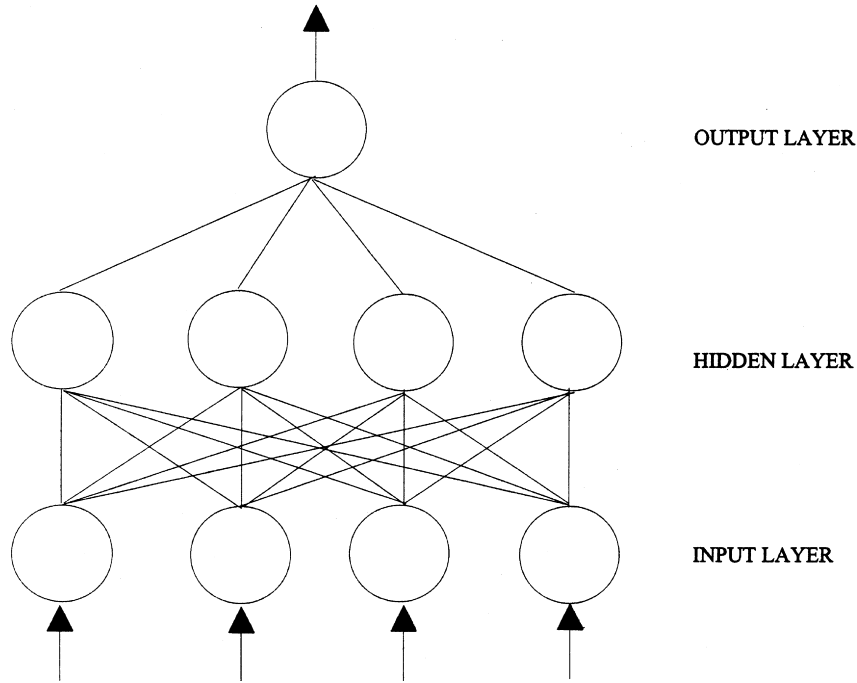


Fig. 1. A typical fully connected feedforward neural network (MLP) used for two-group classification problems.

During the training process, patterns or examples are presented to the input layer of a network. The activation values of the input nodes are weighted and accumulated at each node in the hidden layer. The weighted sum is transferred by an appropriate transfer function into the node's activation value. It then becomes an input into the nodes in the output layer. Finally an output value is obtained to match the desired value. The aim of training is to minimize the differences between the ANN output values and the known target values for all training patterns.

Let $x = (x_1, x_2, \dots, x_n)$ be an n -vector of predictive or attribute variables, y be the output from the network, w_1 and w_2 be the matrices of linking weights from input to hidden layer and from hidden to output layer, respectively. Then a three-layer MLP is in fact a nonlinear model of the form

$$y = f_2(w_2 f_1(w_1 x)), \quad (1)$$

where f_1 and f_2 are the transfer functions for hidden node and output node, respectively. The most popular choice for f_1 and f_2 is the sigmoid function:

$$f_1(x) = f_2(x) = (1 + e^{-x})^{-1}. \quad (2)$$

The purpose of network training is to estimate the weight matrices in Eq. (1) such that an overall error measure such as the mean squared errors (MSE) or sum of squared errors (SSE) is minimized. MSE can be defined as

$$\text{MSE} = \frac{1}{N} \sum_{j=1}^N (a_j - y_j)^2, \quad (3)$$

where a_j and y_j represent the target value and network output for the j th training pattern respectively, and N is the number of training patterns.

From this perspective, network training is an unconstrained nonlinear minimization problem. The most popular algorithm for training is the well-known backpropagation [54] which is basically a gradient steepest descent method with a constant step size. Due to problems of slow convergence and inefficiency with the steepest descent method, many variations of backpropagation have been introduced for training neural networks

[5,13,41]. Recently, Hung and Denton [27] and Subramanian and Hung [59] have proposed to use a general-purpose nonlinear optimizer, GRG2, in training neural networks. The benefits of GRG2 have been reported in the literature for many classification problems [35,42,59]. This study uses a GRG2 based system to train neural networks.

For a two-group classification problem, only one output node is needed. The output values from the neural network (the predicted outputs) are used for classification. For example, a pattern is classified into group 1 if the output value is greater than 0.5, and into group 2 otherwise. It has been shown that the least squares estimate as in the neural networks used in this study yields the posterior probability of the optimal Bayesian classifier [51]. In other words, outputs of neural networks are estimates of the Bayesian posterior probabilities [28]. As will be discussed in the following section, most classification procedures rely on posterior probabilities to classify observations into groups.

2.2. Neural networks and Bayesian classifiers

While neural networks have been successfully applied to many classification problems, the relationship between neural networks and the conventional classification methods is not fully understood in most applications. In this section, we first give a brief overview of the Bayesian classifiers. Then the link between neural networks and Bayesian classifiers is discussed.

Statistical pattern recognition (classification) can be established through Bayesian decision theory [15]. In classification problems, a random pattern or observation $x \in \mathbb{R}^n$ is given and then a decision about its membership is made. Let ω be the state of nature with $\omega = \omega_1$ for group 1 and $\omega = \omega_2$ for group 2. Define

$P(\omega_j)$ = prior probability for an observation x belonging to group j ,

$f(x|\omega_j)$ = conditional probability density function for x given that the pattern belongs to group j ,

where $j = 1, 2$. Using Bayes rule, the posterior probability is

$$P(\omega_j|x) = \frac{f(x|\omega_j)P(\omega_j)}{f(x|\omega_1)P(\omega_1) + f(x|\omega_2)P(\omega_2)}, \quad j = 1, 2. \quad (4)$$

The Bayes decision rule in classification is a criterion such that the overall misclassification error rate is minimized. The misclassification rate for a given x is

$$P(\omega_i|x) = 1 - P(\omega_j|x) \quad \text{if } x \text{ belongs to } \omega_j, \quad i, j = 1, 2.$$

Thus, the Bayesian classification rule can be stated as

Assign x to group k if

$$1 - P(\omega_k|x) = \min_j (1 - P(\omega_j|x))$$

or equivalently

Assign x to group k if

$$P(\omega_k|x) = \max_j P(\omega_j|x). \quad (5)$$

It is now clear that the Bayesian classification rule is based on the posterior probabilities. In the case that $f(x|\omega_j)$ ($j = 1, 2$) are all normal distributions, the above Bayesian classification rule leads to the well-known linear or quadratic discriminant function. See [15] for a detailed discussion.

To see the relationship between neural networks and Bayesian classifiers, we need the following theorem [40].

Theorem 1. Consider the problem of predicting y from x , where x is an n -vector random variable and y is a random variable. The function mapping $F: x \rightarrow y$ which minimizes the squared expected error

$$E[y - F(x)]^2 \quad (6)$$

is the conditional expectation of y given x ,

$$F(x) = E[y|x]. \quad (7)$$

The result stated in the above theorem is the well-known least-squares estimation theory in statistics.

In classification context, if x is the observed attribute vector and y is the true membership value, that is, $y=1$ if $x \in$ group 1; $y = 0$ if $x \in$ group 2, then $F(x)$ becomes

$$\begin{aligned} F(x) &= E[y|x] = 1P(y = 1|x) + 0P(y = 0|x) \\ &= P(y = 1|x) = P(\omega_1|x). \end{aligned} \quad (8)$$

Eq. (8) shows that the least-squares estimate for the mapping function in classification problem is exactly the Bayesian posterior probability.

As mentioned earlier, neural networks are universal function approximators. A neural network in a classification problem can be viewed as a mapping function, $F: \mathbb{R}^n \rightarrow \mathbb{R}$ (see Eq. (2)), where an n -dimensional input x is submitted to the network and a network output y is obtained to make the classification decision. If all the data in the entire population are available for training, then Eqs. (3) and (6) are equivalent and the neural networks produce the exact posterior probabilities in theory. In practice, however, training data are almost always a sample from an unknown population. Thus it is clear that the network output is actually the estimate of posterior probability, i.e. y estimates $P(\omega_1|x)$.

3. Bankruptcy prediction with neural networks

ANNs have been studied extensively as a useful tool in many business applications including bankruptcy prediction. In this section, we present a rather comprehensive review of the literature on the use of ANNs in bankruptcy prediction.

The first attempt to use ANNs to predict bankruptcy is made by Odom and Sharda [38]. In their study, three-layer feedforward networks are used and the results are compared to those of multi-variate discriminant analysis. Using different ratios of bankrupt firms to nonbankrupt firms in training samples, they test the effects of different mixture level on the predictive capability of neural networks and discriminant analysis. Neural networks are found to be more accurate and robust in both training and test results.

Following [38], a number of studies further investigate the use of ANNs in bankruptcy or business failure prediction. For example, Rahimian et al. [49] test the same data set used by Odom and Sharda [38] using three neural network paradigms: backpropagation network, Athena and Perceptron. A number of network training parameters are varied to identify the most efficient training paradigm. The focus of this study is mainly on the improvement in efficiency of the backpropagation algorithm. Coleman et al. [12] also report improved accuracy over that of Odom and Sharda [38] by using their NeuralWare ADSS system.

Salchenberger et al. [55] present an ANN approach to predicting bankruptcy of savings and loan institutions. Neural networks are found to perform as well as or better than logit models across three different lead times of 6, 12 and 18 months. To test the sensitivity of the network to different cutoff values in classification decision, they compare the results for the threshold of 0.5 and 0.2. The information is useful when one expects different costs related to Type I and Type II errors.

Tam and Kiang's paper [61] has had a greater impact on the use of ANNs in general business classification problems as well as in the application of bankruptcy predictions. Based on [60], they provide a detailed analysis of the potentials and limitations of neural network classifiers for business research. Using bank bankruptcy data, they compare neural network models to statistical methods such as linear discriminant analysis, logistic regression, k nearest neighbor and machine learning method of decision tree. Their results show that neural networks are generally more accurate and robust for evaluating bank status.

Wilson and Sharda [66] and Sharda and Wilson [57] propose to use a rigorous experimental design methodology to test ANNs' effectiveness. Three mixture levels of bankrupt and nonbankrupt firms for training set composition with three mixture levels for test set composition yield nine different experimental cells. Within each cell, resampling scheme is employed to generate 20 different pairs of training and test samples. The results more convincingly show the advantages of ANNs rela-

tive to discriminant analysis and other statistical methods.

With a very small sample size (18 bankrupt and 18 nonbankrupt firms), Fletcher and Goss [19] employ an 18-fold cross-validation method for model selection. Although the training effort for building ANNs is much higher, ANNs yield much better model fitting and prediction results than the logistic regression.

In a large scale study, Altman et al. [4] use over 1000 Italian industrial firms to compare the predictive ability of neural network models with that of linear discriminant analysis. Both discriminant analysis and neural networks produce comparable accuracy on holdout samples with discriminant analysis producing slightly better predictions. As discussed in the paper, neural networks have potential capabilities for recognizing the health of companies, but the black-box approach of neural networks needs further studies.

Poddig [44] reports the results from an ongoing study of bankruptcy prediction using two types of neural networks. The MLP networks with three different data preprocessing methods give overall better and more consistent results than those of discriminant analysis. The use of an extension of Kohonen's learning vector quantizer, however, does not show the same promising results as the MLP. Kerling [31], in a related study, compares bankruptcy prediction between France and USA. He reports that there is no significant difference in the correct classification rates for both American and French companies although different accounting rules and financial ratios are employed.

Brockett et al. [10] introduce a neural network model as an early warning system for predicting insurer insolvency. Compared to discriminant analysis and other insurance ratings, neural networks have better predictability and generalizability, which suggests that neural networks can be a useful early warning system for solvency monitoring and prediction.

Boritz et al. [9] use the algorithms of backpropagation and optimal estimation theory in training neural networks. The benchmark models by Altman [2] and Ohlson [39] are employed. Results show that the performance of different classifiers depends on the proportions of bankrupt

firms in the training and testing data sets, the variables used in the models, and the relative cost of Type I and Type II errors. Boritz and Kennedy [8,9] also investigate the effectiveness of several types of neural networks for bankruptcy prediction problems. Different types of ANNs do have varying effects on the levels of Type I and Type II errors. For example, the optimal estimation theory based network has the lowest Type I error level and the highest Type II error level and backpropagation networks have intermediate levels of Type I and II errors while traditional statistical approaches generally have high Type I error and low Type II error levels. They also find that the performance of ANNs is sensitive to the choice of variables and sampling errors.

Kryzanouski and Galler [32] employ the Boltzman machine to evaluate the financial statements of 66 Canadian firms over seven years. Fourteen financial ratios are used in the analysis. The results indicate that the Boltzman machine is an effective tool for neural networks model building. Increasing the training sample size has positive impact on the accuracy of neural networks.

Leshno and Spector [36] evaluate the prediction capability of various ANN models with different data span, neural network architecture and the number of iterations. Their main conclusions are (1) the prediction capability of the model depends on the sample size used for training; (2) different learning techniques have significant effects on both model fitting and test performance; and (3) overfitting problems are associated with large number of iterations.

Lee et al. [34] propose and compare three hybrid neural network models for bankruptcy prediction. These hybrid models combine statistical techniques such as multi-variate discriminant analysis (MDA) and ID3 method with neural networks or combine two different neural networks. Using Korean bankruptcy data, they show that the hybrid systems provide significant better predictions than benchmark models of MDA and ID3 and the hybrid model of unsupervised network and supervised network has the best performance.

Most studies use the backpropagation algorithm [11,38,55,61,64,66] or its variations [43,49] in

training neural networks. It is well known that training algorithms such as the backpropagation have many undesirable features. Piramuthu et al. [43] address the efficiency of network training algorithms. They find that different algorithms do have effects on the performance of ANNs in several risk classification applications. Coats and Fant [11] and Lacher et al. [33] use a training method called “Cascade-Correlation” in a bankruptcy prediction analysis. Compared to MDA or Altman’s *Z* score model, ANNs provide significantly better discriminant ability. Fanning and Cogger [18] compare the performance of a generalized adaptive neural network algorithm (GANNA) and a backpropagation network. They find that GANNA and backpropagation algorithm are comparable in terms of the predictive capability but GANNA saves them time and effort in building an appropriate network structure. Raghupathi [47] conducts an exploratory study to compare eight alternative neural network training algorithms in the domains of bankruptcy prediction. He finds that the Madaline algorithm is the best in terms of correct classifications. However, comparing the Madaline with the discriminant analysis model shows no significant advantage of one over the other. Lenard et al. [35] first apply the generalized reduced gradient (GRG2) optimizer for neural network training in an auditor’s going concern assessment decision model. Using GRG2 trained neural networks results in better performance in terms of classification rates than using backpropagation-based networks.

Based on the pioneering work by Altman [2], most researchers simply use the same set of five predictor variables as in Altman’s original model [11,33,38,49,57,66]. These financial ratios are (1) working capital/total assets; (2) retained earnings/total assets; (3) earnings before interest and taxes/total assets; (4) market value equity/book value of total debt; (5) sales/total assets. Other predictor variables are also employed. For example, Raghupathi et al. [48] use 13 financial ratios previously used successfully in other bankruptcy prediction studies. Salchengerger et al. [55] initially select 29 variables and perform stepwise regression to determine the final five predictors used in neural networks. Tam and Kiang [61] choose 19 financial

variables in their study. Piramuthu et al. [43] use 12 continuous variables and three nominal variables. Alici [1] employs two sets of financial ratios. The first set of 28 ratios is suggested by profile analysis while the second set of nine variables is obtained by using principal component analysis. Boritz and Kennedy [9] test the neural networks with Ohlson’s nine and 11 variables as well as Altman’s five variables. Rudorfer [53] selects five financial ratios from a company’s balance sheet. It is interesting to note that in the literature one study uses as many as 41 independent variables [36] while Fletcher and Goss [19] and Fanning and Cogger [18] use only three variables.

In order to detect maximal difference between bankrupt and nonbankrupt firms, many studies employ matched samples based on some common characteristics in their data collection process. Characteristics used for this purpose include asset or capital size and sales [19,36,63], industry category or economic sector [48], geographic location [55], number of branches, age, and charter status [61]. This sample selection procedure implies that sample mixture ratio of bankrupt to nonbankrupt firms is 50% to 50%.

Most researchers in bankruptcy prediction using neural networks focus on the relative performance of neural networks over other classical statistical techniques. While empirical studies show that ANNs produce better results for many classification or prediction problems, they are not always uniformly superior [46]. Bell et al. [7] report disappointing findings in applying neural networks for predicting commercial bank failures. Boritz and Kennedy [9] have found in their study that ANNs perform reasonably well in predicting business failure but their performance is not in any systematic way superior to conventional statistical techniques such as logit and discriminant analysis. As the authors discussed that there are many factors which can affect the performance of ANNs. Factors in the ANN model building process such as network topology, training method and data transformation are well known. On top of these ANN related factors, other data related factors include the choice of predictor variables, sample size and mixture proportion. It should be pointed out that in most studies, commercial neural net-

work packages are used, which do restrict the users from obtaining a clear understanding of the sensitivity of solutions with respect to initial starting conditions.

4. Design of the study

ANNs are used to study the relationship between the likelihood of bankruptcy and the relevant financial ratios. Two important questions need to be addressed:

- What is the appropriate neural network architecture for a particular data set?
- How robust the neural network performance is in predicting bankruptcy in terms of sampling variability?

For the first question, there are no definite rules to follow since the choice of architecture also depends on the classification objective. For example, if the objective is to classify a given set of objects as well as possible, then a larger network may be desirable. On the other hand, if the network is to be used to predict the classification of unseen objects, then a larger network is not necessarily better. For the second question, we employ a fivefold cross-validation approach to investigate the robustness of the neural networks in bankruptcy prediction. This section will first define variables and the data used in this study. Then a detailed description of the issues in our neural network model building is given. Finally, we illustrate cross-validation methodology used in the study.

4.1. Measures and sample

As described in the previous section, most neural network applications to bankruptcy problems employ the five variables used by Altman [2] and often a few other variables are also injected into the model. This study utilizes a total of six variables. The first five are the same as those in Altman's study – working capital/total assets, retained earnings/total assets, earnings before interest and tax/total assets, market value of equity/total debt, and sales/total assets. The sixth variable, current assets/current liabilities, measures the

ability of a firm in using liquid assets to cover short term obligations. This ratio is believed to have a significant influence on the likelihood of a firm's filing for bankruptcy.

A sample of manufacturing firms that have filed for bankruptcy from 1980 through 1991 is selected from the pool of publicly traded firms in the United States on New York, American and NASDAQ exchanges. These cutoff dates for the 12 year sample period ensure that the provisions of the 1978 Bankruptcy Reform Act have been fully implemented and that the disposition of all bankrupt firms in the sample can be established by the 1994 year end. An extensive search of bankrupt firms is made of the list provided by the Office of the General Counsel of the Security Exchange Commission (SEC) and non-SEC sources such as the Wall Street Journal Index and the Commerce House's Capital Changes Reporter as well as the COMPUSTAT research tapes. Company descriptions and characteristics required for the identification of filing dates are obtained from LEXIS/NEXIS news reports as well as other SEC filings.

The initial search has netted a sample of 396 manufacturing firms that have filed for bankruptcy. The following editing procedures are further implemented to remove sources of confounding in the sample. Firms that (1) have operated in a regulated industry; (2) are foreign based and traded publicly in the US; and (3) have filed bankruptcy previously are excluded from the sample. These sample screenings result in a total of 110 bankrupt manufacturing firms.

In order to highlight the effects of key financial characteristics on the likelihood that a firm may go bankrupt, a matched sample of non-bankrupt firms is selected. Financial information for the three years immediately preceding bankruptcy is obtained from the COMPUSTAT database. Non-bankrupt firms are selected to match with the 110 bankrupt firms in our sample on two key characteristics: two-digit Standard Industrial Classification code and size. Size corresponds to the total assets of a bankrupt firm in the first of the three years before bankruptcy filing. The six financial ratios for the year immediately before the filing of bankruptcy are constructed as independent variables in this study. In summary, we obtained a

matched sample of 220 firms with 110 observations each in the bankrupt and nonbankrupt group.

4.2. Design of neural network model

Currently there are no systematic principles to guide the design of a neural network model for a particular classification problem although heuristic methods such as the pruning algorithm [50], the polynomial time algorithm [52], and the network information technique [65] have been proposed. Since many factors such as hidden layers, hidden nodes, data normalization and training methodology can affect the performance of neural networks, the best network architecture is typically chosen through experiments. In this sense, neural network design is more an art than a science.

ANNs are characterized by their architectures. Network architecture refers to the number of layers, nodes in each layer and the number of arcs. Based on the results from [14,23,37,42], networks with one hidden layer is generally sufficient for

most problems including classification. All networks used in this study will have one hidden layer. For classification problems, the number of input nodes is the number of predictor variables which can be specified by the particular application. For example, in our bankruptcy prediction model, the networks will have six input nodes in the first layer corresponding to six predictor variables. Node biases will be used in the output nodes and logistic activation function will be specified in the networks. In order to attain greater flexibility in modeling a variety of functional forms, direct connections from the input layer to the output layer will be added (see Fig. 2).

The number of hidden nodes is not easy to determine a priori. Although there are several rules of thumb suggested for determining the number of hidden nodes, such as using $n/2$, n , $n + 1$ and $2n + 1$ where n is the number of input nodes, none of them works well for all situations. Determining the appropriate number of hidden nodes usually involves lengthy experimentation since this parameter is problem and/or data dependent. Huang

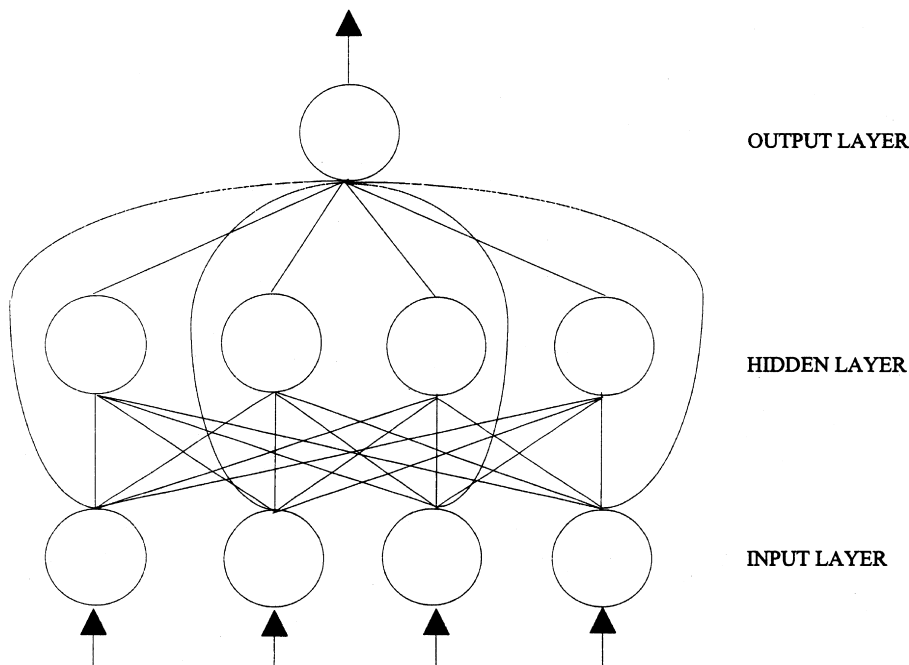


Fig. 2. A complete connected neural network used in this study (direct link from input nodes to the output node).

and Lippmann [26] point out that the number of hidden nodes to use depends on the complexity of the problem at hand. More hidden nodes are called for in complex problems. The issue of the number of hidden nodes also depends on the objective of classification. If the objective is to classify a given set of observations in the training sample as well as possible, a larger network may be desirable. On the other hand, if the network is used to predict classification of unseen objects in the test sample, then a larger network is not necessarily appropriate [42]. To see the effect of hidden nodes on the performance of neural network classifiers, we use 15 different levels of hidden nodes ranging from 1 to 15 in this study.

Another issue in neural networks is the scaling of the variables before training. This so-called data preprocessing is claimed by some authors to be beneficial for the training of the network. Based on our experience (Shanker et al. [56] and also a preliminary study for this project), data transformation is not very helpful for the classification task. Raw data are hence used without any data manipulation.

As discussed earlier, neural network training is essentially a nonlinear nonconvex minimization problem and mathematically speaking, global solutions cannot be guaranteed. Although our GRG2 based training system is more efficient than the backpropagation algorithm [27], it cannot completely eliminate the possibility of encountering local minima. To decrease the likelihood of being trapped in bad local minima, we train each neural network 50 times by using 50 sets of randomly selected initial weights and the best solution of weights among the 50 runs is retained for a particular network architecture.

4.3. Cross-validation

The cross-validation methodology is employed to examine the neural network performance in bankruptcy prediction in terms of sampling variation. Cross-validation is a useful statistical technique to determine the robustness of a model. One simple use of the cross-validation idea is consisted of randomly splitting a sample into two subsam-

ples of training and test sets. The training sample is used for model fitting and/or parameter estimation and the predictive effectiveness of the fitted model is evaluated using the test sample. Because the best model is tailored to fit *one* subsample, it often estimates the true error rate overly optimistically [17]. This problem can be eased by using the so-called fivefold cross-validation, that is, carrying out the simple cross-validation five times. A good introduction to ideas and methods of cross-validation can be found in [20,58].

Two cross-validation schemes will be implemented. First, as in most neural networks classification problems, arc weights from the training sample will be applied to patterns in the test sample. In this study, a fivefold cross-validation is used. We split the total sample into five equal and mutually exclusive portions. Training will be conducted on any four of the five portions. Testing will then be performed on the remaining part. As a result, five overlapping training samples are constructed and testing is also performed five times. The average test classification rate over all five partitions is a good indicator for the out-of-sample performance of a classifier. Second, to have a better picture of the predictive capability of the classifier for the unknown population, we also test each case using the whole data set. The idea behind this scheme is that the total sample should be more representative of the population than a small test set which is only one fifth of the whole data set. In addition, when the whole data set is employed as the test sample, sampling variation in the testing environment is completely eliminated since the same sample is tested five different times. The variability across five test results reflects only the effect of training samples.

The results from neural networks will be compared to those of logistic regression. We choose this technique because it has been shown that the logistic regression is often preferred over discriminant analysis in practice [22,45]. Furthermore, the statistical property of logistic regression is well understood. We would like to know which method gives better estimates of the posterior probabilities and hence leads to better classification results. Since logistic regression is a special case of the neural network without hidden nodes, it is

expected in theory that ANNs will produce more accurate estimates than logistic regression particularly in the training sample. Logistic regression is implemented using SAS procedure LOGISTIC.

5. Results

Table 1 gives the results for the effect of hidden nodes on overall classification performance for both training and small test sets across five subsamples. In general, as expected, one can see when the number of hidden nodes increases, the overall classification rate in the training sets increases. This shows the neural network powerful capability of approximating any function as more hidden nodes are used. However, as more hidden nodes are added, the neural network becomes more complex which may cause the network to learn noises or idiosyncrasies in addition to the underlying rules or patterns. This is recognized as the notorious model overfitting or overspecification problem [21]. For neural networks, obtaining a model that fits the training sample very well is relatively easy if we increase the complexity of a network by, for example, increasing the number of hidden nodes. However, such a large network may have poor generalization capability, that is, it responds incorrectly to other patterns not used in the training process. It is not easy to know a priori when overfitting occurs. One practical way to see this is through the test samples. From Table 1, the best predictive results in test samples are not necessarily those with the larger number of hidden nodes. In fact, neural classifiers with nine or 10 hidden nodes produce the highest classification rates in test samples except for subsample 4 where the best test performance is achieved at four hidden nodes.

For small test sets, cross-validation results on the predictive performance for both neural network models and logistic regression are given in Table 2. This table shows that the overall classification rates of neural networks are consistently higher than those of logistic regression. In addition, neural networks seem to be as robust as logistic regression in predicting the overall classification rate. Across the five small test subsamples,

overall classification rate of neural networks ranges from 77.27% to 84.09% while logistic regression yields classification rates ranging from 75% to 81.82%. However, for each category of bankruptcy and nonbankruptcy, the results indicate no clear patterns. For some subsamples, neural networks predict much better than logistic regression. For others, logistic regression is better. Table 3 gives the pairwise comparison for these two methods in prediction performance. Overall, neural networks are better than logistic regression and the difference of 2.28% is statistically significant at 5% level (p -value is 0.0342). For bankruptcy prediction, neural networks give an average of 81.82% over the five subsamples, higher than 78.18% achieved by logistic regression. For nonbankruptcy prediction, average neural network classification rate is 76.09%, lower than average logistic regression classification rate of 78.18%. Paired t -test results show that the difference between ANNs and logistic regression is not significant in the prediction of bankrupt and nonbankrupt firms.

Tables 4 and 5 show the superiority of ANNs over logistic regression in estimating the true classification rate for the large test set. As we have indicated previously, the large test set is basically the available whole sample data which is consisted of a small test sample and a training sample. Hence, the correct classification rates in Table 4 for the large test set are derived directly from the results for both small test sample and training sample. For example, for training sample 1, the total number of correctly classified firms in the large test set is 191 which is equal to the best small test result (35) plus the corresponding training result (156).

For large test set, ANNs provide consistently not only higher overall classification rates but also higher classification rates for each category of bankrupt and nonbankrupt firms across five training samples. Furthermore, ANNs are more robust than logistic regression in estimating the overall classification rate across five training samples. This is evidenced from the overall classification rate of 86.82% for each of the subsamples 1, 2 and 5, 87.73% for subsample 3, and 85% for subsample 4. Results of paired t -test in Table 5

Table 1
The effect of hidden nodes on overall classification results for training and small test sets ^a

Hidden nodes	Subsample 2			Subsample 3			Subsample 4			Subsample 5		
	Training ^b	Test ^c		Training	Test		Training	Test		Training	Test	
1	139 (78.98)	31 (70.46)		143 (81.25)	36 (81.82)		142 (80.68)	37 (84.09)		142 (80.68)	31 (70.46)	
2	154 (87.50)	28 (63.64)		143 (81.25)	36 (81.82)		142 (80.68)	35 (79.55)		145 (82.39)	34 (77.27)	
3	150 (85.23)	27 (61.36)		146 (82.96)	33 (75.00)		142 (80.68)	37 (84.09)		147 (83.52)	34 (77.27)	
4	148 (84.09)	26 (59.09)		144 (81.82)	27 (61.36)		147 (83.52)	35 (79.55)		151 (85.80)	34 (77.27)	
5	147 (83.52)	29 (65.91)		145 (82.39)	31 (70.46)		146 (82.96)	36 (81.82)		152 (86.36)	35 (79.55)	
6	154 (87.50)	31 (70.46)		154 (87.50)	32 (72.73)		153 (86.93)	37 (84.09)		147 (83.52)	32 (72.73)	
7	156 (88.64)	33 (75.00)		155 (88.07)	35 (79.55)		152 (86.36)	37 (84.09)		154 (87.50)	33 (75.00)	
8	156 (88.64)	29 (65.91)		156 (88.64)	30 (68.18)		153 (86.93)	36 (81.82)		154 (87.50)	33 (75.00)	
9	156 (88.64)	35 (79.55)		155 (88.07)	36 (81.82)		156 (88.64)	35 (79.55)		155 (88.07)	28 (63.64)	
10	158 (89.77)	26 (59.09)		157 (89.21)	31 (70.46)		156 (88.64)	37 (84.09)		156 (88.64)	32 (72.73)	
11	159 (90.34)	24 (54.55)		158 (89.77)	35 (79.55)		157 (89.21)	35 (79.55)		157 (89.21)	34 (77.27)	
12	159 (90.34)	27 (61.36)		159 (90.34)	35 (79.55)		156 (88.64)	37 (84.09)		156 (88.64)	29 (65.91)	
13	159 (90.34)	23 (52.27)		159 (90.34)	26 (59.09)		158 (89.77)	34 (77.27)		157 (89.21)	30 (68.18)	
14	161 (91.48)	26 (59.09)		159 (90.34)	34 (77.27)		158 (89.77)	34 (77.27)		159 (90.34)	32 (72.73)	
15	160 (90.91)	29 (65.91)		160 (90.91)	32 (72.73)		157 (89.21)	33 (75.00)		157 (89.21)	31 (70.46)	
							159 (90.34)	33 (75.00)		160 (90.91)	28 (63.64)	

^a The number in the table is the number of correctly classified; percentage is given in bracket.

^b Training sample size is 176.

^c Test sample size is 44.

Table 2
Cross-validation results on the predictive performance for small test set ^a

Method	Subsample 1 ^b			Subsample 2			Subsample 3			Subsample 4			Subsample 5		
	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall
Neural network	15	20	35	20	16	36	20	17	37	18	17	35	17	17	34
Logistic regression	18	16	34	17	17	34	17	19	36	18	17	35	16	17	33
	(81.82)	(72.73)	(77.27)	(77.27)	(77.27)	(77.27)	(77.27)	(86.36)	(81.82)	(81.82)	(77.27)	(79.55)	(72.73)	(77.27)	(75.00)

^a The number in the table is the number of correctly classified; percentage is given in bracket.

^b B stands for bankruptcy group; NB stands for nonbankruptcy group.

Table 3
Pairwise comparison between ANNs and logistic regression for small test set

Statistics	Overall		Bankrupt		Nonbankrupt	
	ANN	Logistic	ANN	Logistic	ANN	Logistic
Mean	80.46	78.18	81.82	78.18	76.09	78.18
<i>t</i> -statistic	3.1609		0.7182		0.1963	
<i>p</i> -value	0.0342		0.5124		0.8539	

clearly show that the differences between ANNs and logistic regression in the overall and individual class classification rates are statistically significant at the 0.05 level. The differences in overall, bankruptcy and nonbankruptcy classification rate are 8.36%, 11.27% and 4.91%, respectively.

Comparing the results for small test sets in Table 2 and those for large test sets in Table 4, we make the following two observations. First, the variability in results across the five large test samples is much smaller than that of the small test set. This is to be expected as we pointed out earlier that the large test set is the same for each of the five different training sets and the variability in the test results reflects only the difference in the training set. Second, the performance of logistic regression models is stable, while the neural network performance improves significantly, from small test sets to large test sets. The explanation lies in the fact that neural networks have much better classification rates in the training samples. Tables 6 and 7 list the training results of neural networks and logistic regression. The training results for neural networks are selected according to the best overall classification rate in the small test set. Neural networks perform consistently and significantly better for all cases. The differences between ANNs and logistic regression in overall, bankruptcy and nonbankruptcy classification are 9.54%, 13.18% and 5.90%, respectively.

6. Summary and conclusions

Bankruptcy prediction is a class of interesting and important problems. A better understanding of the causes will have tremendous financial and managerial consequences. We have presented a

general framework for understanding the role of neural networks for this problem. While traditional statistical methods work well for some situations, they may fail miserably when the statistical assumptions are not met. ANNs are a promising alternative tool that should be given much consideration when solving real problems like bankruptcy prediction.

The application of neural networks has been reported in many recent studies of bankruptcy prediction. However, the mechanism of neural networks in predicting bankruptcy or in general classification is not well understood. Without a clear understanding of how neural networks operate, it will be difficult to reap full potentials of this technique. This paper attempts to bridge the gap between the theoretical development and the real world applications of ANNs.

It has already been theoretically established that outputs from neural networks are estimates of posterior probabilities. Posterior probabilities are important not only for traditional statistical decision theory but also for many managerial decision problems. Although there are many estimation procedures for posterior probabilities, ANNs is the only known method which estimates posterior probabilities directly when the underlying group population distributions are unknown. Based on the results in this study and [28], neural networks with their flexible nonlinear modeling capability do provide more accurate estimates, leading to higher classification rates than other traditional statistical methods. The impact of the number of hidden nodes and other factors in neural network design on the estimation of posterior probabilities is a fruitful area for further research.

This study used a cross-validation technique to evaluate the robustness of neural classifiers with respect to sampling variation. Model robustness has important managerial implications particularly when the model is used for prediction purposes. A useful model is the one which is robust across different samples or time periods. The cross-validation technique provides decision makers with a simple method for examining predictive validity. Two schemes of fivefold cross-validation methodology are employed. Results show that neural networks are in general quite robust. It is encour-

Table 4
Cross-validation results on the estimation of true classification rates for large test set ^a

Method	Subsample 1 ^b			Subsample 2			Subsample 3			Subsample 4			Subsample 5		
	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall
Neural network	95 (86.36)	96 (87.27)	191 (86.82)	98 (89.09)	93 (84.55)	191 (86.82)	102 (92.73)	91 (82.73)	193 (87.73)	94 (85.45)	93 (84.55)	187 (85.00)	97 (88.18)	94 (85.45)	191 (86.82)
Logistic regression	87 (79.09)	86 (78.18)	173 (78.64)	87 (79.09)	89 (80.91)	176 (80.00)	83 (75.45)	89 (80.91)	172 (78.18)	86 (78.18)	88 (80.00)	174 (79.09)	81 (73.64)	88 (80.00)	169 (76.82)

^a The number in the table is the number of correctly classified; percentage is given in bracket.

^b B stands for bankruptcy group; NB stands for nonbankruptcy group.

Table 5
Pairwise comparison between ANNs and logistic regression for large test set

Statistics	Overall			Bankrupt			Nonbankrupt		
	ANN	Logistic	ANN	ANN	Logistic	ANN	ANN	Logistic	ANN
Mean	86.64	78.55	88.36	88.36	77.09	84.91	84.91	80.00	84.91
t-statistic	10.3807					5.6211		4.0737	
p-value	0.0005					0.0049		0.0152	

Table 6
Comparison of ANNs vs. logistic regression on training sample ^a

Method	Subsample 1 ^b			Subsample 2			Subsample 3			Subsample 4			Subsample 5		
	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall	B	NB	Overall
Neural network	80 (90.91)	76 (86.36)	156 (88.64)	78 (88.64)	77 (87.50)	155 (88.07)	82 (93.18)	74 (84.09)	156 (88.64)	76 (86.36)	76 (86.36)	152 (86.36)	80 (90.91)	77 (87.50)	157 (89.20)
Logistic regression	69 (78.41)	70 (79.55)	139 (78.98)	70 (79.55)	72 (81.82)	142 (80.68)	66 (75.00)	70 (79.55)	136 (77.27)	68 (77.27)	71 (80.68)	139 (78.98)	65 (73.86)	71 (80.68)	136 (77.27)

^a The number in the table is the number of correctly classified; percentage is given in bracket.

^b B stands for bankruptcy group; NB stands for nonbankruptcy group.

Table 7

Pairwise comparison between ANNs and logistic regression for training sample

Statistics	Overall		Bankrupt		Nonbankrupt	
	ANN	Logistic	ANN	Logistic	ANN	Logistic
Mean	88.18	78.64	90.00	76.82	86.36	80.46
<i>t</i> -statistic	9.9623		6.8578		13.8807	
<i>p</i> -value	0.0006		0.0024		0.0002	

aging to note that the variation across samples in training and test classification rates are reasonably small. Much of the variation in results is associated with the number of hidden nodes and initial starting seeds. Users of ANNs will be well advised to use a large number of sets of random starting seeds and experiment on the hidden nodes. After the “optimal” solution is identified and the appropriate number of hidden nodes is selected, the neural classifiers tend to provide consistent estimates.

We also compared neural networks with logistic regression, a well-known statistical method for classification. Neural networks provide significantly better estimate of the classification rate for the unknown population as well as for the unseen part of the population. It can be easily argued that the cost of not being able to predict a bankruptcy is much higher than that for a nonbankrupt firm. Neural networks in our study clearly show their superiority over logistic regression in the prediction of bankrupt firms.

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