ON PATTERN CLASSIFICATION AND SYSTEM IDENTIFICATION BY PROBABILISTIC NEURAL NETWORKS

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In the paper probabilistic neural networks are discussed in detail. Neural network structures for non-parametric pattern classification and system identification are developed.

1. Introduction

If the process is characterized by the absence of a priori information, the most popular methodology for identification and pattern classification is based on non-parametric approach. Such techniques — derived from non-parametric estimates of probability density and regression functions — have been developed by many authors to classify and identify different types of systems (see e.g. Gałkowski and Rutkowski, 1985; 1986; Rutkowski, 1988; 1991; 1993; Rutkowski and Rafajlowicz, 1989). The purpose of this article is to propose neural network structures for implementation of non-parametric algorithms. We shall define a neural network structure as a collection of parallel processors connected together in the form of a directed graph, so organized that the network structure corresponds to the non-parametric problem being considered. We shall use the so-called probabilistic neural networks (Specht, 1990). It should be emphasized that our neural networks do not require learning phase and are asymptotically optimal.

2. Density Estimation

Let $X_1, ..., X_n$ be a sequence of independent, identically distributed random variables taking values in \mathbb{R}^d and having a probability density function f. The Parzen-Rosenblatt estimate of f is given by the formula

$$\widehat{f}_n(x) = \frac{1}{nh_n^d} \sum_{i=1}^n K\left(\frac{x - X_i}{h_n}\right)$$
 (1)

where K is an appriopriately choosen function fulfilling the conditions:

$$\sup_{y} |K(y)| < \infty \tag{2}$$

$$\int_{\mathbb{R}^d} |K(y)| \mathrm{d}y < \infty \tag{3}$$

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$$\lim_{\|\boldsymbol{y}\|\to\infty} \|\boldsymbol{y}\|^d |K(\boldsymbol{y})| = 0 \tag{4}$$

$$\int_{\mathbb{R}^d} K(\boldsymbol{y}) \mathrm{d}\boldsymbol{y} = 1 \tag{5}$$

In eqn. (4) symbol $\|\cdot\|$ stands for the Euclidean vector norm. Sequence h_n is a function of n and satisfies the conditions

$$\lim_{n \to \infty} h_n = 0 \quad \text{and} \quad \lim_{n \to \infty} n h_n^d = \infty$$
 (6)

We assume that the function K is of the form

$$K(\mathbf{x}) = (2\pi)^{-\frac{1}{2}d} e^{-\frac{1}{2}||\mathbf{x}||^2} \tag{7}$$

where $||x||^2 = x^T x$. Then we can rewrite estimator (1) as follows:

$$\widehat{f}_n(x) = \frac{1}{(2\pi)^{d/2} n h_n^d} \sum_{i=1}^n \exp\left(-\frac{(x - X_i)^T (x - X_i)}{2h_n^2}\right)$$
(8)

Observe that

$$(x - X_i)^T (x - X_i) = -2 \left(x^{(1)} X_i^{(1)} + x^{(2)} X_i^{(2)} + \dots + x^{(d)} X_i^{(d)} \right)$$

$$+ \left(x^{(1)} \right)^2 + \left(x^{(2)} \right)^2 + \dots + \left(x^{(d)} \right)^2$$

$$+ \left(X_i^{(1)} \right)^2 + \left(X_i^{(2)} \right)^2 + \dots + \left(X_i^{(d)} \right)^2$$

$$(9)$$

Now assuming normalization of the vectors x and Xi formula (8) simplifies to

$$\widehat{f}_n(x) = \frac{1}{(2\pi)^{d/2} n h_n^d} \sum_{i=1}^n \exp\left(-\frac{(1 - x^T X_i)}{h_n^2}\right)$$
(10)

Figure 1 shows a neural realization of algorithm (10). The proposed net has d inputs and two layers. The first layer consists of n neurons and each neuron has d weights. The output layer has a single neuron with the linear activation function. We should emphasize that the proposed network does not require a training procedure (optimal choosing of connection weights). The succeeding coordinates of the observation vectors X_i , i = 1, ..., n play the role of the weights.

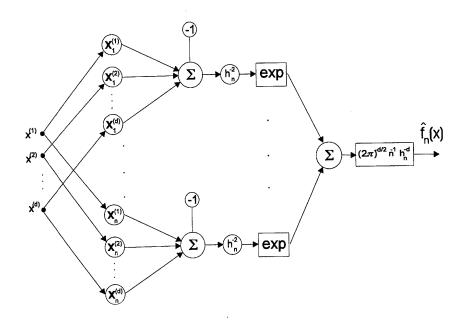


Fig. 1. Neural network density estimation.

The following theorem is a consequence of results presented by Parzen (1962) and Cacoullos (1966).

Theorem 1. If the number of neurons in the first layer of the prabability neural network presented in Fig. 1 is chosen to satisfy conditions (6), then

$$E\left[\widehat{f}_n(\boldsymbol{x}) - f_n(\boldsymbol{x})\right]^2 \xrightarrow{n} 0 \tag{11}$$

in the points at which f is continuous.

3. Pattern Classification

Let (X,Y), (X_1,Y_1) , (X_2,Y_2) , ..., be a sequence of i.i.d. pairs of random variables; Y takes the values from the set; $S = \{1,...,M\}$, whereas X takes the values in \mathbb{R}^d . The problem is to estimate Y from X and V_n , where $V_n = (X_1,Y_1),...,(X_n,Y_n)$ is a learning sequence. Suppose that p_m and f_m , m = 1,...,M, are the prior class probabilities and the class conditional densities, respectively. Define

$$T_{im} = \begin{cases} 1 & \text{if} \quad \mathbf{Y}_i = m \\ 0 & \text{if} \quad \mathbf{Y}_i \neq m \end{cases}$$

for i = 1, 2, ..., n and m = 1, 2, ..., M. The Bayes discriminate function is given by

$$g_m(\mathbf{x}) = f(\mathbf{x})E[T_{nm}|\mathbf{X}_n = \mathbf{x}]$$
(12)

where $f(\mathbf{x}) = \sum_{m=1}^{M} p_m f_m(\mathbf{x})$.

We consider a procedure of classifying every x to a class $m, m \in S$, which maximizes $\widehat{g}_{nm}(x)$, where $\widehat{g}_{nm}(x)$ is the following estimate of the Bayes discriminate function

$$\widehat{g}_{nm}(x) = \frac{1}{(2\pi)^{d/2} n h_n^d} \sum_{i=1}^n T_{im} \exp\left(-\frac{1 - x^T X_i)}{h_n^2}\right)$$
(13)

Figure 2 shows the neural network classifying pattern $x \in \mathbb{R}^d$ to the class m, $m \in S = \{1, 2\}$, for which the expression

$$\widetilde{g}_{nm}(\boldsymbol{x}) = \sum_{i=1}^{n} T_{im} \exp\left(-\frac{1 - \boldsymbol{x}^{T} \boldsymbol{X}_{i})}{h_{n}^{2}}\right)$$

takes the maximum value.

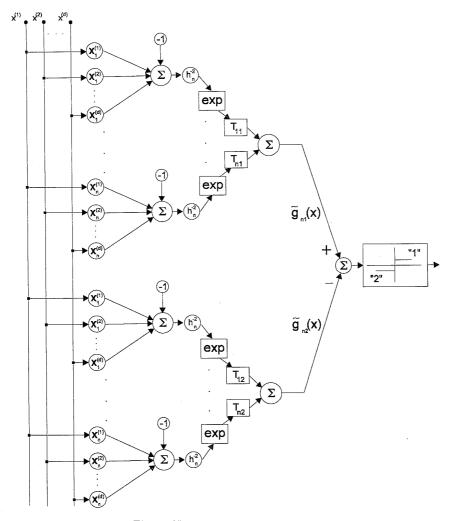


Fig. 2. Neural network pattern classification.

Let $\widehat{\mathbf{Y}}_n$ be a decision obtained according to the classification procedure defined above. Let $D_n = P(\widehat{\mathbf{Y}}_n \neq \mathbf{Y} | \mathbf{V}_n)$. The procedure is said to be weakly (strongly) Bayes risk consistent if $D_n \xrightarrow{n} D_0$ in probability (with probability one), where D_0 is the Bayes probability error. The result below follows from the theorem presented by Greblicki *et al.* (1984).

Theorem 2. If the number of neurons of the probabilistic neural network presented in Fig. 2 satisfies conditions (6), then $D_n \xrightarrow{n} D_0$ in probability.

4. Neural Realization of the Recurrent Non-parametric Estimation Algorithms

The following recurrent estimate of the probability density function was first proposed by Wolverton and Wagner (1969)

$$\widehat{f}_n(\mathbf{x}) = \frac{1}{n} \sum_{i=1}^n \frac{1}{h_i^d} K\left(\frac{\mathbf{x} - \mathbf{X}_i}{h_i}\right)$$
(14)

Expression (14) can be rewritten as

$$\widehat{f}_n(\boldsymbol{x}) = \frac{n-1}{n} f_{n-1}(\boldsymbol{x}) + \frac{1}{nh_n^d} K\left(\frac{\boldsymbol{x} - \boldsymbol{X}_n}{h_n}\right)$$
(15)

For the kernel K of form (7), using the same arguments as those in Section 2, one gets

$$\widehat{f}_n(x) = \frac{n-1}{n} f_{n-1}(x) + \frac{1}{(2\pi)^{d/2} n h_n^d} \exp\left(-\frac{1 - x^T X_n}{h_n^2}\right)$$
(16)

Figure 3 shows the neural network performing algorithm (16) in a fixed point $x \in \mathbb{R}^d$. The net consists of one neuron in the first layer having d inputs — coordinates of the vector X_n , $n = 1, 2, \ldots$. Let us notice that the role of weights is played by the coordinates of the vector x. The second layer also consists of only one neuron with the feedback typical for recurrent neural networks. When the density is to be estimated at several points x_1, \ldots, x_L , then the proposed structure should be copied L times. We shall obtain the neural network of L neurons processing the input observations in the parallel way.

Observe that the recurrent version of algorithm (13) takes the form

$$\widehat{g}_{n,m}(x) = \frac{n-1}{n} \widehat{g}_{n-1,m}(x) + \frac{1}{(2\pi)^{d/2} n h_n^d} T_{n,m} \exp\left(-\frac{1-x^T X_n}{h_n^2}\right)$$
(17)

The corresponding recurrent neural network for pattern classification is shown in Fig. 4.

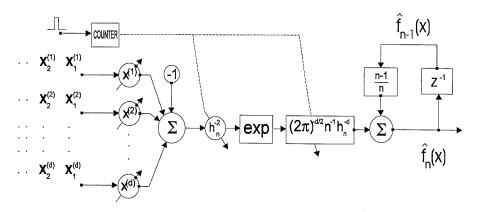


Fig. 3. Neural network density estimation — recurrent algorithm.

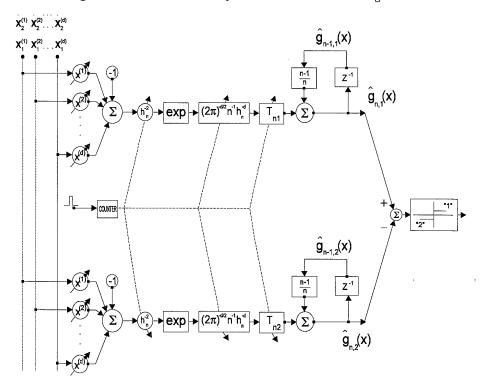


Fig. 4. Neural network pattern classification — reccurent algorithm.

5. Identification by Regression Function Estimation

5.1. Stochastic Input Signal

Let (X, Y) be a pair of random variables. X takes values in \mathbb{R}^d , whereas Y takes values in \mathbb{R} . Let f be the marginal Lebesgue density of X. Based on the sample

 $(X_1, Y_1), ..., (X_n, Y_n)$ of independent observations of (X, Y), we wish to estimate the regression IR of Y on X, i.e. IR(x) = E[Y|X=x]. The probabilistic neural network that provides estimates of function IR and converges to the underlying (linear or non-linear) regression surface is presented in Fig. 4. In system identification the independent variable X is the system input and the dependent variable Y is the system output. Assume that k(x, y) represents the joint probability density function of a vector random variable X and a scalar random variable Y. The conditional mean of Y given x (particular measured value of random variable X) is given by

$$IR(\boldsymbol{x}) = E[\boldsymbol{Y}|\boldsymbol{X} = \boldsymbol{x}] = \frac{1}{f(\boldsymbol{x})} \int_{-\infty}^{+\infty} yk(\boldsymbol{x}, y) \,dy$$
 (18)

If the probability density functions f and k are unknown, then they should be estimated from a sample of observations of X and Y. The joint probability density estimator derived from the Gaussian kernel takes the form:

$$\widehat{k}_n(\boldsymbol{x}, y) = \frac{1}{(2\pi)^{(d+1)/2} n h_n^{d+1}} \sum_{i=1}^n \exp\left(-\frac{(1 - \boldsymbol{x}^T \boldsymbol{X}_i)}{h_n^2}\right) \exp\left(-\frac{(y - \boldsymbol{Y}_i)^2}{2h_n^2}\right)$$
(19)

Replacing k(x, y) by $\hat{k}_n(x, y)$ and f(x) by $\hat{f}_n(x)$ in (18) and performing the indicated integration we get the following estimate of the regression function:

$$\widehat{\mathbf{IR}}_{n}(\boldsymbol{x}) = \frac{\sum_{i=1}^{n} \boldsymbol{Y}_{i} \exp\left(\frac{1 - \boldsymbol{x}^{T} \boldsymbol{X}_{i}}{h_{n}^{2}}\right)}{\sum_{i=1}^{n} \exp\left(\frac{1 - \boldsymbol{x}^{T} \boldsymbol{X}_{i}}{h_{n}^{2}}\right)}$$
(20)

Figure 5 shows a neural network realization of algorithm (20).

One may easily derive the neural network structure for recurrent regression function estimation. The corresponding net is shown in Fig. 6.

Theorem 3. If $E|Y| < \infty$ and the number of neurons of the probabilistic neural network presented in Fig. 5 satisfies conditions (6), then $\widehat{\mathbb{R}}_n \stackrel{n}{\longrightarrow} \mathbb{R}$ in probability for almost all $x \in \mathbb{R}^d$.

This result follows the theorem given by Greblicki et al. (1984).

5.2. Deterministic Input Signal

Many engineering problems are concerned with systems described by the following equation:

$$y_i = IR(x_i) + Z_i, \qquad i = 1, ..., n$$

relating the input x_i and output y_i , and the measurement noise Z_i .

Consider the d-dimensional unit cube $Q = [0, 1]^d$. Let $n^{1/d}$ be an integer, and $i_j = 1, ..., n^{1/d}$, j = 1, ..., d. Partition the interval [0, 1] on each axis into $n^{1/d}$ subsets $\Delta x_{j,i,j}$. Define the following Cartesian product

$$\Delta x_{1,i_1} \times \Delta x_{2,i_2} \times \ldots \times \Delta x_{d,i_d} = Q_{d,i_1}$$

Let $Q_{d,i} \cap Q_{d,j} = \emptyset$, for $i \neq j$ and $\bigcup Q_{d,i} = Q_d$. The inputs x_i are selected so that $x_i \in Q_{d,i}$.

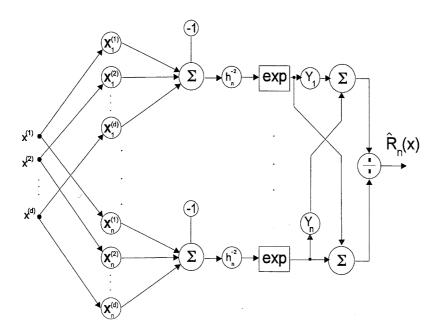


Fig. 5. Neural regression function estimation — stochastic input signal.

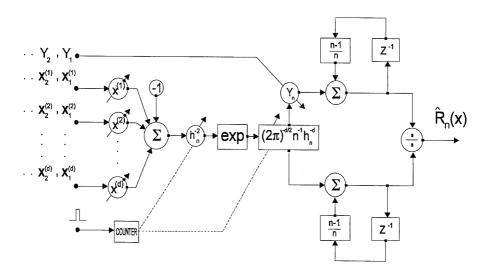


Fig. 6. Recurrent neural regression estimation.

We propose the following algorithm

$$\widehat{\mathbf{R}}_n(\boldsymbol{x}) = \frac{1}{(2\pi)^{d/2} n h_n^d} \sum_{i=1}^n \boldsymbol{Y}_i \exp\left(-\frac{(1-\boldsymbol{x}^T \boldsymbol{X}_i)}{h_n^2}\right)$$
(21)

The appropriate net is shown in Fig. 7.

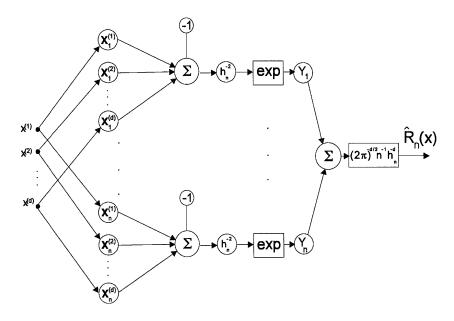


Fig. 7. Neural regression function estimation — deterministic input signal.

Theorem 4. If the number of neurons of the probabilistic neural network presented in Fig. 7 satisfies conditions (6) and $\sup_{1 \le i \le n} \sup_{x,y \in Q_i} ||x-y|| = O(n^{-1/d})$, then $\widehat{\mathbb{R}}_n \xrightarrow{n} \mathbb{R}$ in probability at all continuity points x of \mathbb{R} .

The above theorem follows from theorems presented by Gałkowski and Rutkowski (1985) and Georgiev (1990).

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