

# A LEARNING APPROACH TO IDENTIFICATION OF NONLINEAR PHYSIOLOGICAL SYSTEMS USING WIENER MODELS

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Abstract: The Wiener model is a natural description of many physiological systems. Although there have been a number of algorithms proposed for the identification of Wiener models, most of the existing approaches were developed under some restrictive assumptions of the system such as a white noise input, part or full invertibility of the nonlinearity, or known nonlinearity. In this study a new recursive algorithm based on Lyapunov stability theory is presented for the identification of Wiener systems with unknown and noninvertible nonlinearity and noisy data. The new algorithm can guarantee global convergence of the estimation error to a small region around zero and is as easy to implement as the well-known back propagation algorithm. Theoretical analysis and example studies show the effectiveness and advantages of the proposed method compared with the earlier approaches.

## 1 INTRODUCTION

Numerous approaches have been proposed for the identification of nonlinear systems including parametric and nonparametric methods (Greblicki 1997, Nelles 2001). Among these, the so-called block-oriented models have been found very useful in practice, due to their simplicity in structure and relative ease of implementation and interpretation. One of the block-oriented structures is known as the Wiener model, which consists of a cascade connection of a linear time invariant (LTI) system followed by a static (memoryless) nonlinearity. Such a structure has been shown to be a reasonable model for many chemical and biological processes (e.g.: Hunter and Korenberg 1986), as well as communication and control systems (Huang 1998, Bloemen et al 2001). Theoretically, any nonlinear system that has a Volterra or Wiener functional expansion can be represented (with a sufficient degree of accuracy) by a finite sum of Wiener

models (Boyd and Chua 1985).

Several different algorithms have been presented in the literature for the identification of Wiener models. Early approaches used correlation analysis, but long periods of data and white Gaussian noise inputs are required (Billings and Fakhouri 1978). Approaches based on the invertibility of the static nonlinearity, and estimation of the linear and nonlinear blocks either in a successive (Narendra et al 1966) and iterative procedure or in a simultaneous manner (Gomez and Baeyens 2004, Kalafatis 1997) have also been proposed. The main disadvantage of such algorithms is that convergence is difficult to guarantee. Moreover, several studies assumed the nonlinearity to be known (Wigren 1994) or approximated by a piecewise linear function (the nonlinearity needs to be invertible in each of the small working regions identified - Figueroa 2008). Similarly, Bai and Reyland (2009) assumed the nonlinearity to be monotonic (and therefore invertible) in a specific region. Only a few studies do not assume and make use of the invertibility of

the nonlinear block. Lacy and Bernstein (2003) directly expanded the system into a “linear in parameters” regressive form. The approach also requires additional manipulation to extract the model parameters for the linear and nonlinear part. Comparisons made by the authors with previous approaches showed that their singular value decomposition SVD-based method and the gradient-based algorithm provide better estimates. Nonetheless, the algorithms are computationally expensive, especially when the orders of the linear and nonlinear parts are high.

In the present contribution a learning approach for the identification of Wiener models with unknown and non-invertible nonlinearity, based on Lyapunov stability theory is proposed. Previous work has studied the identification of nonlinear systems using learning methods based on neural networks (Kosmatopoulos et al 1995). However the use of the learning approach for direct identification of Wiener models from input-output data has not been fully explored. The proposed recursive algorithm is developed with guaranteed global convergence. The linear part is given by an IIR or FIR filter model and the nonlinear part is approximated by a polynomial. All model parameters are estimated simultaneously, and linear and nonlinear model orders can be set to be arbitrarily high. The new approach is as simple as a back propagation (BP) algorithm with regard to implementation. The learning approach can also be used to estimate time-varying systems, which is of particular relevance to the neurophysiological investigations that motivated the current work. Theoretical analysis and simulation results to evaluate the effectiveness of the method are also presented.

## 2 WIENER MODEL IDENTIFICATION PROBLEM

The Wiener model is composed of a linear block followed by a static nonlinear unit (Fig.1). The linear part is assumed to be single-input single-output (SISO) linear IIR model. The Wiener system can therefore be written as:

$$x(t) = -a_1x(t-1) - a_2x(t-2), \dots - a_{N_a}x(t-N_a) \quad (1)$$

$$+ b_0u(t) + b_1u(t-1) + \dots + b_{N_b}u(t-N_b) \quad (2)$$

$$y(t) = f(x(t)) + w(t)$$

where  $u(t)$ ,  $x(t)$  and  $y(t)$  are the input to the system, the (unmeasured) output of the linear part, and the

measured output of the system, respectively. In the process, input and output noise can all be regarded as additive output noise denoted by  $w(t)$ . The nonlinear function is assumed to be a polynomial function of the form:

$$f(x) = c_0 + c_1x + c_2x^2 + \dots + c_{N_c}x^{N_c} \quad (3)$$

Note that a polynomial function with sufficiently high order can be used to approximate any continuous nonlinearity to any degree of accuracy in the region of interest for  $x$  (Jeffreys 1988). Here, the nonlinearity  $f(\cdot)$  is not necessarily invertible.

For convenience, (1-3) can be written as:

$$\begin{aligned} x(t) &= K^T U_t, & f(x(t)) &= C^T X_t, \\ y(t) &= f(K^T U_t) + w(t) \end{aligned} \quad (4)$$

$$\begin{aligned} \text{where } K &= [a_1, a_2, \dots, a_{N_a}, b_0, b_1, \dots, b_{N_b}]^T \\ U_t &= [-x(t-1), \dots, -x(t-N_a), u(t), u(t-1), \dots, u(t-N_b)]^T \\ C &= [c_0, c_1, c_2, \dots, c_{N_c}]^T, & X_t &= [1, x, x^2, \dots, x^{N_c}] \end{aligned} \quad (5)$$

with  $N_a$ ,  $N_b$  and  $N_c$  the corresponding orders used in estimation. The estimation error can be defined as

$$e(t) = \hat{y}(t) - y(t) = \hat{f}(\hat{x}(t)) - f(x(t)) - w(t) \quad (6)$$

The identification problem is to find an updated law for the model in (4-5)

$$\hat{K}(t) = \hat{K}(t-1) + \Delta \hat{K}(t) \quad (7a)$$

$$\hat{C}(t) = \hat{C}(t-1) + \Delta \hat{C}(t) \quad (7b)$$

given a series of input-output data pairs  $u(t)$  and  $y(t)$  ( $t=1, 2, \dots, T$ ), with any initial values  $\hat{K}(0)$  and  $\hat{C}(0)$ , such that the estimation error in (6) comes to zero (noise-free case) or a small region near zero (noisy case) as  $t \rightarrow \infty$ , according to a cost function  $V(e(t))$  which is a positive definite function of  $e(t)$ . Thus assuming stationary signals and a time-invariant system, each model parameter converges to a constant level. To ensure a unique solution,  $\hat{K}(t)$  and  $\hat{C}(t)$  can be normalized. For example if the linear part is estimated as an FIR model and suppose  $\hat{b}_0 \neq 0$ :

$$\begin{aligned} \hat{K}(t) &= \hat{K}(t) / \hat{b}_0 \\ \hat{C}(t) &= [\hat{c}_0, \hat{b}_0 \hat{c}_1, \hat{b}_0^2 \hat{c}_2, \dots, \hat{b}_0^{N_c} \hat{c}_{N_c}]^T \end{aligned} \quad (8)$$

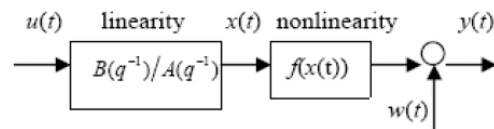


Figure 1: Wiener model.

### 3 THE LEARNING METHOD

The learning method (LM) updates the model parameters with each new sample, driving a cost function towards zero. The algorithm, based on Lyapunov stability theory, is formulated as follows:

**Lemma 1.** The difference of the estimation error (6) between two successive sampling times can be computed as

$$\Delta e(t) = e(t) - e(t-1) = \Delta \hat{y}(t) - \Delta y(t) \quad (9)$$

where  $\Delta$  indicates the change between successive samples and  $\hat{\cdot}$  an estimate. By expanding  $\hat{f}(\cdot)$  as a Taylor series:

$$\begin{aligned} \Delta e(t) &= \hat{X}_i^T \Delta \hat{C}(t) + \hat{f}_i(t) \Delta \hat{K}(t) + \varepsilon(t) + \sigma(\cdot) \\ \text{where } \hat{f}_i(t) &= \frac{\partial \hat{f}(t)}{\partial \hat{x}(t)} = \hat{C}(t-1)^T \frac{\partial \hat{X}_i(t)}{\partial \hat{x}(t)} \\ \hat{f}_{ii}(t) &= \frac{\partial^2 \hat{f}(t)}{\partial \hat{x}(t)^2} = \hat{C}(t-1)^T \frac{\partial^2 \hat{X}_i(t)}{\partial \hat{x}(t)^2} \\ \varepsilon(t) &= \hat{f}_i(t) \hat{K}(t-1)^T \Delta \hat{U}_i \\ &+ \frac{1}{2} \hat{f}_{ii}(t) \Delta \hat{U}_i^T \hat{K}(t-1)^T \Delta \hat{U}_i - y(t) + y(t-1) \end{aligned} \quad (10)$$

where  $\sigma(\cdot) = \sigma(\Delta \hat{C}(t), \Delta \hat{K}(t), \Delta \hat{U}_i)$  denotes the remaining higher order terms in a Taylor series expansion of  $\hat{y}(t)$  and measurement noise, and  $\Delta \hat{U}_i = \hat{U}_i - \hat{U}_{i-1}$ .

**Remark 1.** In (10) the effects of model parameter updates (first two terms) and effect of the changing input (represented by  $\varepsilon(t)$ ) on errors are explicitly considered. Note that the conventional back propagation (BP) algorithm in learning methods is simply based on the assumption that the output error has no distinct relationship with the input  $u(t)$ , therefore limiting the use of BP for the identification of Wiener models. The current method thus overcomes this important limitation of a conventional approach.

**Theorem 1.** Given input output data pairs  $u(t)$  and  $y(t)$  ( $t=1,2,\dots, n \gg \max(N_u, N_y)$ ) measured for system (4) and with the assumption that  $|\sigma(\cdot)| < \rho$ , the estimated model (4 and 5) can be obtained with the estimation error (6) asymptotically convergent to a ball with radius  $\sqrt{\rho \sigma_a / \eta_c}$ , ( $\eta_c > 0, \sigma_a > 0$ ) around zero by training the estimation model with the parameter update laws (7a,b).

Proof of this theorem will be presented elsewhere.

**Remark 2.** The new algorithm is globally convergent, in terms of a cost function  $V(t) = e^2(t)$ , to a small region around zero whose size is determined by the upper bound of  $\sigma(t)$  which

denotes the remaining higher order terms in a Taylor series expansion and also represents the “effect” of the model estimation error. Existing recursive two-step methods (i.e. Hunter and Korenberg 1986) can not guarantee convergence and the recursive algorithm in Wigren (1993) can only guarantee it locally. It should also be emphasized that the algorithm proposed does not require the nonlinearity to be invertible

**Remark 3.** When there is additive noise in the measured output, the error (6) will not represent the true difference in output between the real and the estimated model. This will affect the update laws in (7a,b) and thus result in  $\sigma(t)$ , due to the high order terms of the Taylor series, to vary with a larger amplitude ( $\rho$ ). Setting  $\eta_k$  (the learning rate for the linear parameters in  $\hat{K}$ ) as small as possible will reduce the problem. Note that the convergence speed of the algorithm is mainly determined by  $\eta_c$  (the learning rate for the nonlinear parameters in  $\hat{C}$ ). Also, the saturation-like error  $\bar{e}(t)$  is used to avoid the unnecessary oscillations in the recursive computation which might arise following sudden large errors.

$$\begin{aligned} \Delta \hat{K}(t) &= \begin{cases} -\eta_k \frac{1}{\hat{f}_i(\cdot)} \frac{\hat{U}_i}{\hat{U}_i^T \hat{U}_i} \bar{e}(t) & \text{if } |\hat{f}_i(\cdot)| > \delta \\ \eta_k \operatorname{sgn}(\eta_k) & \text{otherwise} \end{cases} \\ \Delta \hat{C}(t) &= \begin{cases} -(\eta_c - \eta_k) \frac{\hat{X}_i}{\hat{X}_i^T \hat{X}_i} \bar{e}(t) \\ -\frac{\hat{X}_i}{\hat{X}_i^T \hat{X}_i} (\bar{e}(t) - \gamma(t)) & \text{if } |\hat{f}_i(\cdot)| > \delta \\ -(\eta_c - \eta_k |\hat{f}_i(\cdot)|) \frac{\hat{X}_i}{\hat{X}_i^T \hat{X}_i} \bar{e}(t) \\ -\frac{\hat{X}_i}{\hat{X}_i^T \hat{X}_i} (\bar{e}(t) - \gamma(t)) & \text{otherwise} \end{cases} \quad (11) \\ \gamma(t) &= -\rho \frac{e(t)}{\max\{|e(t)|, \sigma_a\}}, \quad \operatorname{sgn}(e) = \begin{cases} 1 & \text{if } |e| > 0 \\ 0 & \text{if } |e| = 0 \\ -1 & \text{if } |e| < 0 \end{cases} \end{aligned}$$

$$\bar{e}(t) = \begin{cases} e_b \operatorname{sgn}(e(t)) & \text{if } |e(t)| > e_a \\ e(t) & \text{otherwise} \end{cases}, \quad e_b \geq e_a \gg \sigma_a > 0$$

$$\bar{\varepsilon}(t) = \begin{cases} \varepsilon_a \operatorname{sgn}(\varepsilon(t)) & \text{if } |\varepsilon(t)| > \varepsilon_a \\ \varepsilon(t) & \text{otherwise} \end{cases}, \quad \varepsilon_b > 0$$

### 4 EXAMPLES

**Example 1.** Consider a Wiener model  $K=[a_1, b_0, b_1]^T=[1, 1, 2]^T$  with a noninvertible nonlinearity,  $C=[c_0, c_1, c_2, c_3, c_4]^T=[0.0001 \ 0.0010 \ 0.0150 \ -0.0005 \ -0.0001]^T$  as shown in Fig. 2. The system is stimulated by white Gaussian noise with added white noise such that the signal to noise ratio is 2, i.e.,  $\|y_r\|/\|w\|=2$ . The table in Figure 2a shows

the model parameters estimated with our proposed learning method LM after three training rounds (which are sufficient for the algorithm to converge provided the learning rate is appropriately selected). These are compared with the SVD method (Lacy 2003). The results show that even though slightly larger errors are obtained for  $b_1$  and  $c_0$  using the proposed algorithm, all other parameter errors are considerably better than those by the SVD method (see Fig 2a). The model fit for the validation data (not used in training) is 61.58% with the LM whilst is only 38.22% with the SVD. The fitness to the real output without noise is 96.41% for the LM and only 47.24% for SVD.

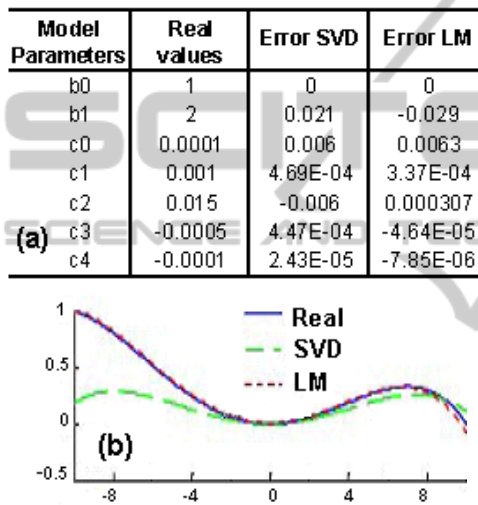


Figure 2: a) Errors in model estimates. b) Nonlinearity to be identified.

**Example 2.** The LM algorithm was also applied to the intracellular potential recorded from a spiking local interneuron, that is part of the reflex control loop of the hind limb (Newland et al 1997, Vidal-Gadea et al 2009). The input signal was Gaussian noise used to stimulate a stretch-sensor located at the femoro-tibial joint of the hind leg. The noninvertible nonlinearity identified using the proposed learning method is shown in Figure 3a. The fitness for validation data is 50.0% after three rounds of training. The LM algorithm was also run in a BP-like condition whereby the consideration of the effect from the changing input (Remarks 1-3) was removed. In this case, the fitness in the same validation data is only 38.67% (three rounds of training). Here the model orders were  $N_a=10$ ,  $N_b=30$  and  $N_c=9$ . Due to the high orders of the model, it is difficult to apply the SVD method.

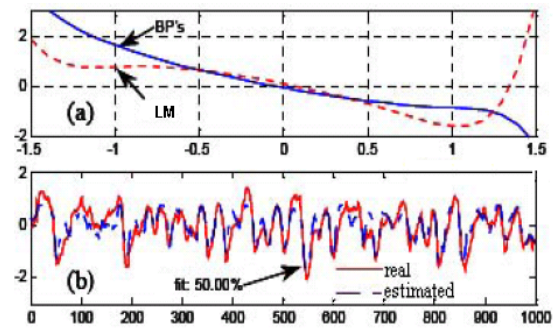


Figure 3: A practical example from a locust neuro muscular control systems. (a) The estimated nonlinearity, (c) Estimated output (LM).

## 5 CONCLUSIONS

Most of the existing algorithms for the identification of Wiener models were developed under some restrictive assumptions, such as white noise input, part or full invertibility of the nonlinearity, or known nonlinearity. A novel recursive algorithm based on a learning approach has been developed for the identification of Wiener systems with unknown and noninvertible nonlinearity and noisy data. The new algorithm can guarantee global convergence of the estimation error to a small range around zero and is easy to implement in a manner similar to the well-known back propagation (BP) algorithm. Comparisons between the proposed methodology and existing algorithms such as SVD-based method and BP algorithm were provided in two example studies. The theoretical analysis and example studies show the effectiveness and advantages of the proposed approach. In continuing this work, we will investigate optimal choices of the control parameters for the algorithm, and provide more extensive evaluations in simulated and recorded signals.

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