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Refined Instrumental Variable Methods for Identifying Hammerstein Models Operating in Closed Loop

V. Laurain*, M. Gilson*, H. Garnier*

Abstract—This article presents an instrumental variable method dedicated to non-linear Hammerstein systems operating in closed loop. The linear process is a Box–Jenkins model and the non-linear part is a sum of known basis functions. The performance of the proposed algorithm is illustrated by a numerical example.

I. INTRODUCTION

The need for non-linear identification grows as the studied system complexity increases [14]. Many different approaches were developed to deal with open-loop black-box model identification, whether they are non parametric, using Volterra series approach [11], [15], semi-parametric using neural network methods and support vector machine classification [16], [3], or parametric such as state dependent parameters [24] or refined instrumental variable [12]. Semi-parametric approaches, even if performing efficiently, lack the possibility of giving an *a posteriori* physical representation of the studied system. On the other hand, transfer function models provide a generic approach to data-based modelling of linear systems, encompass both discrete-time and continuous-time applications and are in an ideal form to interpret serial, parallel connections of sub-systems which often have a physical significance.

Closed-loop systems are widely used in practice and various attempts have been made to handle linear system identification in the presence of feedback. Indeed, closed-loop system identification leads to several difficulties due to the correlation between the disturbances and the control signal induced by the loop. Several methods have therefore been developed to cope with this problem see *e.g.* [18], [20], [5], [25]. An optimal refined instrumental variable method has been presented for closed-loop linear systems in [7].

The problem of non-linear plants operating in closed loop have not received a lot of attention so far. Indeed, this problem is handled only in a few papers as in *e.g.* [6], [10], [2], and [21] but to the best of our knowledge no instrumental variable (IV) method has been developed to deal with this kind of issues. For closed-loop linear identification, a basic IV estimator was first suggested in [22], assuming the knowledge of the controller; the topic was later discussed in more detail in [19]. The present paper extends the refined IV technique theory developed for linear cases presented in [7] to the identification of discrete-time Hammerstein models operating in closed loop. This paper is organized as follows: after the preliminaries, the proposed iterative IV method

Fig. 1. Closed-loop system

is described in Section III. The detailed algorithm of the method is given in Section IV. Finally, the performance of the proposed algorithm is illustrated through a numerical example in Section V.

II. PROBLEM DESCRIPTION

A. System considered

Consider the stable, SISO closed-loop non-linear system represented in Figure 1 and assume that both command input and output signals, $u(t_k)$ and $y(t_k)$ are uniformly sampled at a constant sampling time T_s over N samples.

The controller is supposed to be known, the linear process is Box-Jenkins model and the non-linear function is assumed to be a sum of known basis functions $\gamma_1, \gamma_2, \dots, \gamma_l$ so that it can be expressed:

$$\bar{u}(t_k) = \sum_{i=1}^l \alpha_{o,i} \gamma_i(u(t_k)). \quad (1)$$

Notice first that this system produces the same input-output data for any pair $(\beta f(u), G_o(q)/\beta)$. Therefore, to get a unique parametrization, one of the gains of $f(u)$ or $G_o(q)$ has to be fixed [4], [1]. Hence, the first coefficient of the function $f(\cdot)$ equals to 1, *i.e.* $\alpha_1 = 1$ in (1).

Moreover, $\bar{u}(t_k)$ in (1) is an internal variable and is not directly available from data. Using (1), the system $\mathcal{S}$ to be identified is described by the following input-output relationships:

$$\mathcal{S} \begin{cases} y(t_k) = G_o(q)\bar{u}(t_k) + H_o(q)e_o(t_k), \\ \bar{u}(t_k) = \sum_{i=1}^l \alpha_{o,i} \gamma_i(u(t_k)), \\ u(t_k) = r(t_k) - C_c(q)y(t_k), \end{cases} \quad (2)$$

where $G_o(q)$ denotes the process and $C_c(q)$ the controller. $u(t_k)$ describes the process input signal, $y(t_k)$ the process output signal. $B_o(q^{-1})$ and $A_o(q^{-1})$ are polynomials in the shift operator q^{-1} ($q^{-r}x(t_k) = x(t_{k-r})$) with respective degree n_b and n_a . This non-linear system can be expressed

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in the following form by combining the first two equations in (2):

$$\mathcal{S} \begin{cases} y(t_k) = \sum_{i=1}^l G_{o,i}(q) \gamma_i(u(t_k)) + H_o(q) e_o(t_k), \\ u(t_k) = r(t_k) - C_c(q) y(t_k), \end{cases} \quad (3)$$

where

$$G_{o,i}(q) = \frac{B_{o,i}(q^{-1})}{A_o(q^{-1})} = \frac{\alpha_{o,i} B_o(q^{-1})}{A_o(q^{-1})}, \quad (4)$$

and

$$r(t_k) = r_1(t_k) + C_c(q) r_2(t_k). \quad (5)$$

The method presented is based on the identification of a Box–Jenkins model, where the linear and the noise models are not constrained to have common polynomials. The coloured noise associated with the sampled output measurement $y(t_k)$ has rational spectral density and can be represented by a discrete-time autoregressive moving average ARMA model:

$$\xi(t_k) = H_o(q) e(t_k) = \frac{C_o(q^{-1})}{D_o(q^{-1})} e_o(t_k) \quad (6)$$

where $C_o(q^{-1})$ and $D_o(q^{-1})$ are polynomials with respective degree n_c and n_d . $e_o(t_k)$ is a zero-mean, normally distributed, discrete-time white noise sequence: $e_o(t_k) \sim \mathcal{N}(0, \sigma_e^2)$.

Additionally, the controller $C_c(q)$ is given by

$$C_c(q) = \frac{Q(q^{-1})}{P(q^{-1})} = \frac{q_0 + q_1 q^{-1} + \dots + q_{n_q} q^{-n_q}}{p_0 + p_1 q^{-1} + \dots + p_{n_p} q^{-n_p}}, \quad (7)$$

with the pair (P, Q) assumed to be coprime.

B. Model considered

The model set to be estimated, denoted as $\mathcal{M}$ with subsystems $(\mathcal{G}_i)$ and noise $(\mathcal{H})$ models parameterized independently, then takes the form,

$$\mathcal{M} : \{G_i(q, \boldsymbol{\rho}), H(q, \boldsymbol{\eta})\}, \quad \text{with } i = 1 \dots l \quad (8)$$

where $\boldsymbol{\rho}$ and $\boldsymbol{\eta}$ are parameter vectors that characterise the system and noise model, respectively. In particular, the system model is formulated in discrete-time terms:

$$\begin{aligned} \mathcal{G} : G_i(q, \boldsymbol{\rho}) &= \frac{B_i(q^{-1}, \boldsymbol{\rho})}{A(q^{-1}, \boldsymbol{\rho})}, \\ &= \frac{\alpha_i (b_1 q^{-1} + b_2 q^{-2} + \dots + b_{n_b} q^{-n_b})}{1 + a_1 q^{-1} + \dots + a_{n_a} q^{-n_a}}, \end{aligned} \quad (9)$$

with $i = 1 \dots l$ and the pairs (A, B_i) assumed to be coprime. The associated model parameters are stacked columnwise in the parameter vector,

$$\boldsymbol{\rho} = \begin{bmatrix} \mathbf{a} \\ \alpha_1 \mathbf{b} \\ \vdots \\ \alpha_l \mathbf{b} \end{bmatrix} \in \mathbb{R}^{n_\rho}, \mathbf{a} = \begin{bmatrix} a_1 \\ a_2 \\ \vdots \\ a_{n_a} \end{bmatrix} \in \mathbb{R}^{n_a}, \mathbf{b} = \begin{bmatrix} b_1 \\ b_2 \\ \vdots \\ b_{n_b} \end{bmatrix} \in \mathbb{R}^{n_b}, \quad (10)$$

with $n_\rho = n_a + n_b l$ while the noise model is in discrete-time form

$$\mathcal{H} : H(q, \boldsymbol{\eta}) = \frac{C(q^{-1}, \boldsymbol{\eta})}{D(q^{-1}, \boldsymbol{\eta})} = \frac{1 + c_1 q^{-1} + \dots + c_{n_c} q^{-n_c}}{1 + d_1 q^{-1} + \dots + d_{n_d} q^{-n_d}} \quad (11)$$

where the associated model parameters are stacked columnwise in the parameter vector,

$$\boldsymbol{\eta} = [c_1 \quad \dots \quad c_{n_c} \quad d_1 \quad \dots \quad d_{n_d}]^T \in \mathbb{R}^{n_c + n_d} \quad (12)$$

Consequently, the noise transfer function takes the usual ARMA model form:

$$\xi(t_k) = \frac{C(q^{-1}, \boldsymbol{\eta})}{D(q^{-1}, \boldsymbol{\eta})} e(t_k). \quad (13)$$

More formally, there exists the following decomposition of the parameter vector $\boldsymbol{\theta}$ for the whole model,

$$\boldsymbol{\theta} = \begin{pmatrix} \boldsymbol{\rho} \\ \boldsymbol{\eta} \end{pmatrix} \quad (14)$$

such that the model can be written in the form

$$\begin{cases} y(t_k) = \frac{1}{A(q^{-1}, \boldsymbol{\rho})} \sum_{i=1}^l B_i(q^{-1}, \boldsymbol{\rho}) \gamma_i(u(t_k)) + \xi(t_k) \\ \xi(t_k) = \frac{C(q^{-1}, \boldsymbol{\eta})}{D(q^{-1}, \boldsymbol{\eta})} e(t_k), \\ u(t_k) = r(t_k) - C_c(q) y(t_k) \end{cases} \quad (15)$$

with $B_i(q^{-1}, \boldsymbol{\rho}) = \alpha_i B(q^{-1}, \boldsymbol{\rho})$, $i = 1 \dots l$.

C. Identification problem

Under the following assumptions:

- A1: The degrees n_a , n_b , n_c and n_d are supposed to be known *a priori*;
- A2: The controller C_c is supposed to be *a priori* known and will not be estimated;
- A3: $A_o(q^{-1})$ and $B_o(q^{-1})$ are coprime;
- A4: $r(t_k)$ is not correlated to $e_o(t_k)$;
- A5: $r(t_k)$ and $y(t_k)$ are stationary signals;
- A6: The non-linear basic functions γ_i as well as the degree l are *a priori* known and γ_i are pairwise orthogonal;

The identification problem can be formulated as follows: given N samples of data $\{r(t_k), u(t_k), y(t_k)\}_{k=1}^N$, directly estimate the discrete time parameter vector $\boldsymbol{\theta}$.

III. INSTRUMENTAL VARIABLE FOR CLOSED-LOOP SYSTEMS

A. The instrumental variable estimate

Consider the relationship between process input and output signals which can be written under such a regression form:

$$y(t_k) = \boldsymbol{\varphi}^T(t_k) \boldsymbol{\rho} + v(t_k) \quad (16)$$

The well-known extended-IV estimate is given by [17]:

$$\hat{\rho}_{xiv}(N) = \arg \min_{\boldsymbol{\rho}} \left\| \left[\frac{1}{N} \sum_{k=1}^N L(q) \zeta(t_k) L(q) \boldsymbol{\varphi}^T(t_k) \right] - \left[\frac{1}{N} \sum_{k=1}^N L(q) \zeta(t_k) L(q) y(t_k) \right] \right\|_W^2, \quad (17)$$

where $\zeta(t_k)$ is the instrumental vector, $\|x\|_W^2 = x^T W x$, with W a positive definite weighting matrix and $L(q)$ a stable prefilter. By definition, when $G_o \in \mathcal{G}$, the extended-IV estimate provides a consistent estimate under the following two conditions¹

- $\mathbb{E} L(q) \zeta(t_k) L(q) \boldsymbol{\varphi}^T(t_k)$ is full column rank,
- $\mathbb{E} L(q) \zeta(t_k) L(q) v(t_k) = 0$.

B. Optimal instrumental variable for linear systems in closed loop

Concerning the latest advance in IV method for closed-loop systems, a so-called “tailor-made IV algorithm” was proposed in [8], where the closed-loop plant is parameterized using the (open-loop) plant parameters. Then, an optimal (minimal) variance result was developed in the linear closed-loop extended IV identification case, revealing consequences for the choice of weights, filters and instruments [9], [7].

Consider the relationship between the process input and output signals in (2) for $u(t) = \bar{u}(t)$,

$$y(t_k) = G_o(q) u(t_k) + H_o(q) e_o(t_k). \quad (18)$$

If the process to be identified belongs to the model set previously defined, then $y(t_k)$ can be written under the regression form:

$$y(t_k) = \boldsymbol{\varphi}^T(t_k) \boldsymbol{\rho} + v(t_k) \quad (19)$$

where,

$$\boldsymbol{\varphi}(t_k) = [-y(t_{k-1}) \dots -y(t_{k-n_a}), u(t_{k-1}) \dots u(t_{k-n_b})]^T, \quad (20)$$

$$\boldsymbol{\rho} = [a_1, \dots, a_{n_a}, b_1, \dots, b_{n_b}]^T \quad (21)$$

and

$$v(t_k) = A(q^{-1}, \boldsymbol{\rho}) \xi(t_k). \quad (22)$$

It has been shown in that case that the minimum variance estimator can be achieved by the following optimal choice of design variables [9], [7]

$$W = I, \quad (23)$$

$$\zeta(t_k) = \dot{\varphi}(t_k), \quad (24)$$

$$L(q) = L^{opt}(q) = \frac{1}{A_o(q^{-1}) H_o(q)}. \quad (25)$$

where $\dot{\varphi}(t_k)$ is the noise-free version of $\varphi(t_k)$.

¹The notation $\mathbb{E}[\cdot] = \lim_{N \rightarrow \infty} \frac{1}{N} \sum_{t=1}^N \mathbb{E}[\cdot]$ is adopted from the prediction error framework of [13].

C. Instrumental variable for non-linear systems in closed loop

Now consider the relationship between the process input and output signals in (2),

$$y(t_k) = G_o(q) \bar{u}(t_k) + H_o(q) e_o(t_k). \quad (26)$$

If the process to be identified belongs to the model set previously defined, then $y(t_k)$ can be written under the regression form (with $\boldsymbol{\rho}$ as defined in (10)) :

$$y(t_k) = \boldsymbol{\varphi}^T(t_k) \boldsymbol{\rho} + v(t_k) \quad (27)$$

where,

$$\boldsymbol{\varphi}(t_k) = \begin{bmatrix} -\mathbf{y}(t_k) \\ \gamma_1(\mathbf{u}(t_k)) \\ \vdots \\ \gamma_l(\mathbf{u}(t_k)) \end{bmatrix}, \mathbf{y}(t_k) = \begin{bmatrix} y(t_{k-1}) \\ y(t_{k-2}) \\ \vdots \\ y(t_{k-n_a}) \end{bmatrix}, \quad (28)$$

$$\mathbf{u}(t_k) = \begin{bmatrix} u(t_{k-1}) \\ u(t_{k-2}) \\ \vdots \\ u(t_{k-n_b}) \end{bmatrix}, \quad (29)$$

and

$$v(t_k) = A(q^{-1}, \boldsymbol{\rho}) \xi(t_k). \quad (30)$$

By logically extending the IV presented in Section III-B to non-linear systems, we propose to use the following instrumental variable:

$$\zeta(t_k) = \begin{bmatrix} -\dot{\mathbf{y}}(t_k) \\ \gamma_1(\dot{\mathbf{u}}(t_k)) \\ \vdots \\ \gamma_l(\dot{\mathbf{u}}(t_k)) \end{bmatrix}, \dot{\mathbf{y}}(t_k) = \begin{bmatrix} \dot{y}(t_{k-1}) \\ \dot{y}(t_{k-2}) \\ \vdots \\ \dot{y}(t_{k-n_a}) \end{bmatrix}, \quad (31)$$

$$\dot{\mathbf{u}}(t_k) = \begin{bmatrix} \dot{u}(t_{k-1}) \\ \dot{u}(t_{k-2}) \\ \vdots \\ \dot{u}(t_{k-n_b}) \end{bmatrix},$$

where the noise-free system can be described by:

$$\dot{\mathbf{S}} \begin{cases} \dot{y}(t_k) = \sum_{i=1}^l G_{o,i}(q) \gamma_i(\dot{u}(t_k)), \\ \dot{u}(t_k) = r(t_k) - C_c(q) \dot{y}(t_k) \end{cases} \quad (32)$$

Remark 1

It has to be noticed that for the non-linear case, choosing $L(q)$, $\zeta(t_k)$ and W as optimally defined in the linear case does not guarantee the statistical optimality of the estimate. In the present context it can be seen that n_φ and therefore n_ζ over-parametrizes the system: The number of parameters of the system which equals $n_a + n_b + l - 1$ is less than the number of parameters to be estimated ($n_a + n_b l$) in the associated model.

Remark 2

None of $A(q^{-1}, \boldsymbol{\rho})$, $B_i(q^{-1}, \boldsymbol{\rho})$, $C(q^{-1}, \boldsymbol{\eta})$ or $D(q^{-1}, \boldsymbol{\eta})$ is known and only their estimates are available. Therefore,

refined IV estimation normally involves an iterative (or relaxation) algorithm in which, at each iteration, the ‘auxiliary model’ used to generate the instrumental variable, as well as the associated prefilter $L(q)$, are updated, based on the parameter estimates obtained at the previous iteration [23].

IV. THE REFINED INSTRUMENTAL VARIABLE ALGORITHM

Step 1. Initialisation: A simple least square method is used to initialize the algorithm.

$$\hat{\rho}^0 = \left[\sum_{k=1}^N \varphi(t_k) \varphi^T(t_k) \right]^{-1} \sum_{k=1}^N \varphi(t_k) y(t_k), \quad (33)$$

where $\varphi(t_k)$ is given by (28). This yields $B_i(q^{-1}, \hat{\rho}^0)$ and $A(q^{-1}, \hat{\rho}^0)$. Denote the corresponding transfer functions

$$G_i(q, \hat{\rho}^0) = \frac{B_i(q^{-1}, \hat{\rho}^0)}{A(q^{-1}, \hat{\rho}^0)}. \quad (34)$$

Set the initial noise model estimates $C(q^{-1}, \hat{\eta}^0) = D(q^{-1}, \hat{\eta}^0) = 1$ and $j = 1$.

Step 2. Estimate a model based proposed IV

Generate the filtered instruments according to the model structure at the j^{th} iteration:

- Compute $L(q, \hat{\theta}^{j-1}) = \frac{1}{A(q^{-1}, \hat{\rho}^{j-1})H(q, \hat{\eta}^{j-1})}$,
- Compute the auxiliary model variables $\hat{y}(t_k, \hat{\rho}^{j-1})$ and $\hat{u}(t_k, \hat{\rho}^{j-1})$ by simulating the estimated auxiliary model $\hat{S}(\hat{\rho}^{j-1})$:

$$\begin{cases} \hat{y}(t_k, \hat{\rho}^{j-1}) = \sum_{i=1}^l G_{o,i}(q, \hat{\rho}^{j-1})(\gamma_i(\hat{u}(t_k, \hat{\rho}^{j-1}))), \\ \hat{u}(t_k, \hat{\rho}^{j-1}) = r(t_k) - C_c(q)\hat{y}(t_k, \hat{\rho}^{j-1}), \end{cases} \quad (35)$$

- Generate the filtered instrument:
 $\hat{\zeta}_f(t_k) = L(q, \hat{\theta}^{j-1})\dot{\varphi}(t_k, \hat{\rho}^{j-1})$,
- Generate the filtered regressor and output
 $\varphi_f(t_k) = L(q, \hat{\theta}^{j-1})\varphi(t_k, \hat{\rho}^{j-1})$, and
 $y_f(t_k) = L(q, \hat{\theta}^{j-1})y(t_k)$
- Compute the non-linear closed-loop refined instrumental variable (HCLRIV) estimate:

$$\hat{\rho}^j(N) = \left[\sum_{k=1}^N \hat{\zeta}_f(t_k) \varphi_f^T(t_k) \right]^{-1} \sum_{k=1}^N \hat{\zeta}_f(t_k) y_f(t_k). \quad (36)$$

Step 3. Obtain an estimate of the noise model parameter vector $\hat{\eta}^j$ based on the estimated noise sequence:

- Compute the estimated noise sequence:
 $v(t_k) = y(t_k) - \sum_{i=1}^l \frac{B_i(q^{-1}, \hat{\rho}^j)}{A(q^{-1}, \hat{\rho}^j)} \gamma_i(u(t_k))$,
- Estimate $\hat{\eta}^j$ and therefore $C(q^{-1}, \hat{\eta}^j)$ and $D(q^{-1}, \hat{\eta}^j)$ with an ARMA method using $v(t_k)$ as input. The

ARMA method from the MATLAB identification toolbox is used here.

Step 4. repeat from step 2. Stop when $A(q^{-1}, \hat{\rho})$, $B_i(q^{-1}, \hat{\rho})$, $H(q, \hat{\eta})$ have converged.

Step 5. Compute the estimated α coefficients and $B(q^{-1}, \hat{\rho})$ by using a constraint

At the end of the iterative process, the coefficients $\hat{\alpha}_i$ can be deduced from the polynomial $\hat{B}_i(q^{-1})$ since $B_i(q^{-1}, \rho) = \alpha_i B(q^{-1}, \rho)$. In the present case we use a straightforward constraint, which is $\alpha_1 = 1$ (guarantees that $\hat{B}_1(q^{-1}, \hat{\rho}) = \hat{B}(q^{-1}, \hat{\rho})$) to compute $\hat{\alpha}_i$ from:

$$\hat{\alpha}_i = \frac{1}{n_b + 1} \sum_{k=0}^{n_b} \frac{\hat{b}_{i,k}}{\hat{b}_{1,k}}, \quad (37)$$

where $\hat{b}_{i,k}$ is the k th coefficient of the polynomial term $B_i(q^{-1}, \hat{\rho})$.

Comments

- A simplified version of HCLRIV algorithm named HCLSRIV follows the exact same theory for estimation of Hammerstein models operating in closed-loop with a linear output error process. It is mathematically described by, $C(q^{-1}, \eta^j) = C_o(q^{-1}) = 1$ and $D(q^{-1}, \eta^j) = D_o(q^{-1}) = 1$. All previous steps remain true except for *step 3* which is not anymore required.
- Even if the proposed algorithm performs well, it is not statistically optimal as illustrated in Section V. However, the estimated parameters can be used to initialise a statistically optimal method (prediction error method for example).

V. NUMERICAL EXAMPLE

This section presents numerical illustration of both suggested HCLRIV and HCLSRIV methods. For all presented examples, the non-linear block is a sum of sine functions, i.e. $\gamma_i(u(t_k)) = \sin(i \cdot u(t_k))$, $\forall i \in \mathbb{N}$ and the command input $\bar{u}$ considered here is:

$$\bar{u}(t_k) = \sin(u(t_k)) - 0.5 \sin(2u(t_k)) + 0.4 \sin(3u(t_k)). \quad (38)$$

$r_1(t_k)$ follows a uniform distribution with values between -2 and 2 while $r_2(t_k) = 0$ ($N = 2000$). The linear model and controller assuming a zero order hold on the input are given as:

$$\begin{aligned} G_o(q) &= \frac{0.0997q^{-1} - 0.0902q^{-2}}{1 - 1.8858q^{-1} + 0.9048q^{-2}}, \\ H_o(q) &= \frac{1 + 0.5q^{-1}}{1 - 0.85q^{-1}}, \\ C_c(q) &= \frac{10.75 - 9.25q^{-1}}{1 - q^{-1}}. \end{aligned} \quad (39)$$

The models considered for estimation are:

$$\mathcal{M}_{HCLRIV} \begin{cases} G(q, \rho) = \frac{b_1 q^{-1} + b_2 q^{-2}}{1 + a_1 q^{-1} + a_2 q^{-2}}, \\ H(q, \eta) = \frac{1 + c_1 q^{-1}}{1 + d_1 q^{-1}}, \\ f(u(t_k)) = \sin(u(t_k)) + \sum_{i=2}^3 \alpha_i \sin(i \cdot u(t_k)) \end{cases} \quad (40)$$

for the HCLRIV method and

$$\mathcal{M}_{HCLSRIV} \begin{cases} G(q, \rho) = \frac{b_1 q^{-1} + b_2 q^{-2}}{1 + a_1 q^{-1} + a_2 q^{-2}}, \\ H(q, \eta) = 1, \\ f(u(t_k)) = \sin(u(t_k)) + \sum_{i=2}^3 \alpha_i \sin(i \cdot u(t_k)) \end{cases} \quad (41)$$

for the HCLSRIV method.

As pointed out in Section III the proposed estimator is not optimal due to redundancy in the regressor. For illustration purposes, the proposed methods are compared to the parameters of the linear plant process identified using a prediction error minimisation method (PEMlin, OElin) assuming that α_i are known. The problem reduces therefore to the identification of a linear model where the input is $\bar{u}(t_k)$ and the output is $y(t_k)$.

The model considered for the PEM identification using the PEM method from the MATLAB identification toolbox is then:

$$\mathcal{M}_{PEMlin} \begin{cases} G(q, \rho) = \frac{b_1 q^{-1} + b_2 q^{-2}}{1 + a_1 q^{-1} + a_2 q^{-2}}, \\ H(q, \eta) = \frac{1 + c_1 q^{-1}}{1 + d_1 q^{-1}}, \\ f(u(t_k)) = \sin(u(t_k)) - 0.5 \sin(2u(t_k)) \\ \quad + 0.4 \sin(3u(t_k)) \end{cases} \quad (42)$$

for the PEMlin method.

The model considered for the PEM identification using the OE method from the MATLAB identification toolbox is then:

$$\mathcal{M}_{OElin} \begin{cases} G(q, \rho) = \frac{b_1 q^{-1} + b_2 q^{-2}}{1 + a_1 q^{-1} + a_2 q^{-2}}, \\ H(q, \eta) = 1, \\ f(u(t_k)) = \sin(u(t_k)) - 0.5 \sin(2u(t_k)) \\ \quad + 0.4 \sin(3u(t_k)) \end{cases} \quad (43)$$

for the OElin method.

A Monte-Carlo simulation of 100 runs ($N_{exp} = 100$) with a new noise realization for each run is accomplished under a signal to noise ratio (SNR) of 10dB with:

$$SNR = 10 \log \left(\frac{P_g}{P_e} \right), \quad (44)$$

P_g being the average power of signal g . The number of samples is $N = 2000$. Table I exhibits the mean value of

the estimated parameters, their standard deviation and their normalised root mean square error (RMSE) defined as:

$$RMSE(\hat{\theta}_j) = \sqrt{\frac{1}{N_{exp}} \sum_{i=1}^{N_{exp}} \left(\frac{\theta_j^o - \hat{\theta}_j(i)}{\theta_j^o} \right)^2}, \quad (45)$$

with $\hat{\theta}_j$ the j th component of the parameter vector θ .

On the one hand it can be seen in Table I that even if the noise model assumption is false, which is the case for the HCLSRIV and OElin methods in this example, the refined IV method gives a parameter estimation centered on the true value with a correct estimated parameter variance (RMSE remaining under 16%), while the prediction error based method OElin estimates are strongly biased and have RMSE near 30%. This fact makes HCLSRIV a very interesting method for practical applications where the noise dynamic is not known: by modelling a white noise added to the output, the number of parameters to estimate is reduced in comparison to a Box-Jenkins model identification.

On the other hand, the presented HCLRIV method is not optimal due to the redundancy in the regressor but exposes RMSE in estimated parameters under 11 %. It can be noticed however that the estimated parameter variances are close to those from PEMlin estimates: this may therefore be seen as an acceptable achievement since in the PEMlin case, the non-linear part is supposed to be known.

VI. CONCLUSION

The theory of refined instrumental variable was applied to a non-linear closed-loop model composed of a known linear controller, a linear dynamic Box-Jenkins transfer function and a non-linear function defined as the sum of known basis functions. The proposed estimator is not statistically optimal but it was shown on some given example that the variance of estimated parameters is close to the one obtained from the linear optimal PEM estimator. The proposed method can be therefore used as a good starting point for some statistically optimal PEM methods which are known to be sensitive to the initialisation stage. In case the noise model assumed is false, the proposed IV method still converges while the PEM based methods do not. Some further work will be dedicated to blind estimation problems, where the non-linear basic functions γ_i are *a priori* not known.

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TABLE I
ESTIMATION RESULTS FOR DIFFERENT MODELS

method		b_1	b_2	a_1	a_2	α_1	α_2	c_1	d_1
	true value	0.0997	-0.0902	-1.8858	0.9048	-0.5000	0.4000	0.5000	-0.8500
PEMlin	$mean(\hat{\theta})$	0.0998	-0.0901	-1.8857	0.9028	X	X	0.4918	-0.8603
	$std(\hat{\theta})$	0.0009	0.0025	0.0247	0.0208	X	X	0.0208	0.0312
	RMSE	0.0089	0.0280	0.0131	0.0229	X	X	0.0421	0.0365
OElin	$mean(\hat{\theta})$	0.0693	-0.0653	-1.8655	0.8867	X	X	X	X
	$std(\hat{\theta})$	0.0016	0.0019	0.0186	0.0171	X	X	X	X
	RMSE	0.2959	0.2686	0.0141	0.0267	X	X	X	X
HCLSRIV	$mean(\hat{\theta})$	0.0978	-0.0894	-1.8715	0.8927	-0.4893	0.4006	X	X
	$std(\hat{\theta})$	0.0100	0.0138	0.0943	0.0912	0.0774	0.0495	X	X
	RMSE	0.1010	0.1588	0.0520	0.1042	0.1554	0.1230	X	X
HCLRIV	$mean(\hat{\theta})$	0.1004	-0.0895	-1.8862	0.8923	-0.5024	0.4015	0.4959	-0.8515
	$std(\hat{\theta})$	0.0031	0.0090	0.0829	0.0799	0.0182	0.0153	0.0249	0.0310
	RMSE	0.0313	0.1035	0.0457	0.0915	0.0367	0.0381	0.0496	0.0365

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