

## An optimal instrumental variable method for continuous-time fractional model identification

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**Abstract:** this paper deals with continuous-time system identification using fractional differentiation models in a noisy output context. The simplified refined instrumental variable for continuous-time systems (*srivc*) is extended to fractional models. Monte Carlo simulation analysis are used to demonstrate the performance of the proposed optimal instrumental variable scheme.

Keywords: continuous-time models; fractional differentiation; instrumental variable; state variable filter; system identification; time-domain.

### 1. INTRODUCTION

Instrumental Variable (IV) approaches to system identification and parameter estimation have been developed since 1960s and 1970s as detailed in the survey paper by Young (1981) and all references therein. Recently, Refined Instrumental Variables for Continuous-system (Young and Jakeman (1980)) has been developed for identifying hybrid continuous-time Box-Jenkins models, where the system model is continuous-time and the noise model is discrete-time (Young et al. (2006)). *rivc* algorithms are implemented in both Captain (Young (2006)) and Contsid (Garnier et al. (2008)) toolboxes.

The last two decades have witnessed considerable development in the use of fractional operators in various fields. Fractional differentiation is now an important tool for the international scientific and industrial communities especially in modeling viscoelastic materials and some diffusive phenomena (thermal diffusion, electrochemical diffusion). Time-domain system identification using fractional models was initiated in the late nineties and in the beginning of this century in the Ph.D. theses of Le Lay (1998); Lin (2001); Cois (2002); Aoun (2005). For an overview of different identification methods based on fractional models, refer to Malti et al. (2006).

**The objective of this paper** is to extend the Simplified Refined Instrumental Variable for Continuous-time systems (Young and Jakeman (1980)), denoted by *srivc* from hereon, to the identification of fractional models. This optimal IV method presents the advantage of yielding asymptotically efficient estimates in the presence of white measurement noise. The algorithm is initialized by a transfer function obtained from the least squares state variable filter estimates. Cois et al. (2001) are the first to

extend the use of state variable filters to fractional systems. The authors suggest to use Instrumental Variables (IV) to eliminate bias. However, their instruments are not selected in an optimal way. An improvement of their method is proposed in this paper by choosing *rivc* algorithm. This approach involves a method of iterative-adaptive prefiltering based on a quasi-optimal<sup>1</sup> statistical solution. The proposed method is independent from the way fractional differentiation and integration are simulated in the time-domain.

#### 1.1 Mathematical background

A fractional mathematical model is based on a fractional differential equation:

$$y(t) + a_1 \mathbf{D}^{\alpha_1} y(t) + \dots + a_{m_A} \mathbf{D}^{\alpha_{m_A}} y(t) = b_0 \mathbf{D}^{\beta_0} u(t) + b_1 \mathbf{D}^{\beta_1} u(t) + \dots + b_{m_B} \mathbf{D}^{\beta_{m_B}} u(t), \quad (1)$$

where  $(a_j, b_i) \in \mathbb{R}^2$ , differentiation orders  $\alpha_1 < \alpha_2 < \dots < \alpha_{m_A}$ ,  $\beta_0 < \beta_1 < \dots < \beta_{m_B}$  are allowed to be non-integer positive numbers. The concept of differentiation to an arbitrary order (non-integer),

$$\mathbf{D}^\gamma \triangleq \left( \frac{d}{dt} \right)^\gamma \quad \forall \gamma \in \mathbb{R}_+^*,$$

was defined in the 19<sup>th</sup> century by Riemann and Liouville. The  $\gamma$ -order fractional derivative of  $x(t)$  is defined as being an integer derivative of order  $\lceil \gamma \rceil + 1$  ( $\lceil \cdot \rceil$  stands for the floor operator) of a non-integer integral of order  $\gamma - \lceil \gamma \rceil$  Samko et al. (1993):

<sup>1</sup> As stated in Young et al. (2006), in the rational case, the method is quasi-optimal because true optimality would require optimal interpolation of the input signal  $u(t)$  over the sampling interval, whereas only simple interpolation is used here.

$$\mathbf{D}^\gamma x(t) = \mathbf{D}^{[\gamma]+1} \left( \mathbf{I}^{[\gamma]+1-\gamma} x(t) \right) \triangleq \left( \frac{d}{dt} \right)^{[\gamma]+1} \left( \frac{1}{\Gamma([\gamma]+1-\gamma)} \int_0^t \frac{x(\tau) d\tau}{(t-\tau)^{\gamma-[\gamma]}} \right), \quad (2)$$

where  $t > 0$ ,  $\forall \gamma \in \mathbb{R}_+^*$ , and the Euler's  $\Gamma$  function is defined as:

$$\Gamma(x) = \int_0^\infty e^{-t} t^{x-1} dt \quad \forall x \in \mathbb{R}^* \setminus \{\mathbb{N}^-\}. \quad (3)$$

The Laplace transform is a more concise algebraic tool generally used to represent fractional systems, see Oldham and Spanier (1974):

$$\mathcal{L}\{\mathbf{D}^\gamma x(t)\} = s^\gamma X(s) \quad \text{if } x(t) = 0 \quad \forall t \leq 0. \quad (4)$$

This property allows to write the fractional differential equation (1), provided  $u(t)$  and  $y(t)$  equal 0 for all  $t < 0$ , in a transfer function form:

$$F(s) = \frac{B(s)}{A(s)} = \frac{\sum_{i=0}^{m_B} b_i s^{\beta_i}}{1 + \sum_{j=1}^{m_A} a_j s^{\alpha_j}}. \quad (5)$$

Moreover if  $F(s)$  is commensurable of order  $\gamma$ , i.e. all differentiation orders are exactly divisible by the same number, an integral number of times (the biggest number is always chosen), then  $F(s)$  can be rewritten as:

$$F(s) = \frac{\sum_{i=0}^n \tilde{b}_i s^{i\gamma}}{1 + \sum_{j=1}^m \tilde{a}_j s^{j\gamma}}, \quad (6)$$

where  $n = \frac{\beta_{m_B}}{\gamma}$  and  $m = \frac{\alpha_{m_A}}{\gamma}$  are integers and:

$$\begin{cases} \tilde{b}_i = b_i & \text{if } i\gamma = \beta_i & \text{and } \tilde{b}_i = 0 & \text{if } i\gamma \neq \beta_i \\ \tilde{a}_j = a_j & \text{if } j\gamma = \alpha_j & \text{and } \tilde{a}_j = 0 & \text{if } j\gamma \neq \alpha_j. \end{cases} \quad (7)$$

In rational transfer functions  $\gamma$  equals 1 and usually numerator  $\alpha_{m_A}$  and denominator  $\beta_{m_B}$  orders are both fixed, then all coefficients  $b_i, i = 1, \dots, \beta_{m_B}$  and  $a_j, j = 1, \dots, \alpha_{m_A}$  are estimated. Generally, no care is taken to check whether any intermediate coefficient, as in (7) equals zero.

Exactly the same principle will be applied here for fractional models: the commensurable order  $\gamma$  together with the  $\alpha_{m_A}$  and  $\beta_{m_B}$  are fixed, and then all the  $\tilde{b}_i$  and  $\tilde{a}_j$  coefficients are estimated.

### 1.2 Time-domain simulation of fractional models

The main objective of this paper is to develop a method for time-domain system identification able to work with any fractional-model-time-domain simulation algorithm. Hence, only the simulation algorithm used in the example of this paper is explained here, as the objective is not to discuss different simulation algorithms. Refer to Aoun et al. (2004), for an extended discussion on this topic.

Due to the consideration that real physical systems generally have bandlimited fractional behavior and due to the practical limitations of input and output signals (Shannon's cut-off frequency for the upper band and the spectrum of the input signal for the lower band), fractional

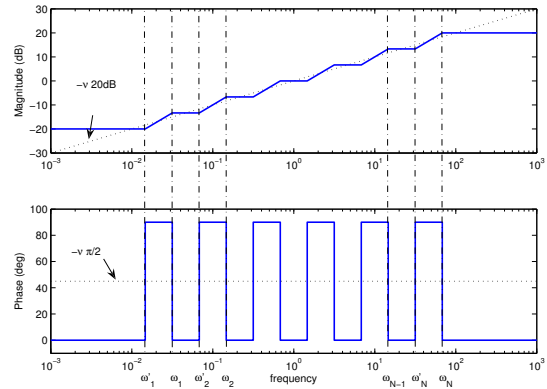


Fig. 1. Recursive approximation of a fractional differentiator with poles and zeroes

operators are usually approximated by high order rational models. As a result, a fractional model and its rational approximation have the same dynamics within a limited frequency band. The most commonly used approximation of  $s^\nu$  in the frequency band  $[\omega_A, \omega_B]$  is the recursive distribution of zeros and poles proposed by Oustaloup (1995) (see Fig. 1):

$$s^\gamma \rightarrow s_{[\omega_A, \omega_B]}^\gamma = C_0 \left( \frac{1 + \frac{s}{\omega_A}}{1 + \frac{s}{\omega_B}} \right)^\gamma \approx C_0 \prod_{k=1}^N \frac{1 + \frac{s}{\omega_k'}}{1 + \frac{s}{\omega_k}} \quad (8)$$

where  $\omega_i = \alpha \omega_i'$ ,  $\omega_{i+1}' = \eta \omega_i'$  and

$$\gamma = 1 - \frac{\log \alpha}{\log \alpha \eta}, \quad (9)$$

$\alpha$  and  $\eta$  define the differentiation order  $\gamma$ . The bigger  $N$  the better the approximation of the differentiator  $s^\nu$ .

## 2. SYSTEM IDENTIFICATION

The **R**efined **I**nstrumental **V**ariables for **C**ontinuous system (Young and Jakeman (1980)) (*rivc*) method and its simplified version *srivc* (when the additive measurement noise is assumed to be white) are the only IV methods that can be interpreted in optimal statistical terms, so providing an estimate of the parametric error covariance matrix and, therefore, estimates of the confidence bounds on the parameter estimates. Optimal instrumental variables are obtained from an auxiliary model which is iteratively updated. At the first iteration, the auxiliary model is initialized using a least squares estimate applied on state variable filters (SVF). The identification problem is first stated, then SVF method is briefly recalled and finally the *srivc* algorithm is extended to fractional models.

### 2.1 Problem formulation

Consider input  $u(t)$  and output  $y^*(t)$  data collected at regular samples  $T_s, 2T_s, \dots, KT_s$  from  $t = 0$  to  $t = T$ . The measured output is supposed to be corrupted by an additive measurement noise  $p(t)$ :

$$y^*(t) = y(t) + p(t), \quad (10)$$

where  $y(t)$  is the hypothetical noise-free deterministic system output. The input  $u(t)$  and the output  $y(t)$  signals are supposed to be related by the fractional differential equation (1).

Equation error can be formulated as:

$$\varepsilon(t) = y^*(t) - \varphi^*(t)^T \theta, \quad (11)$$

where

$$\varphi^*(t)^T = \begin{bmatrix} \mathbf{D}^{\beta_0} u(t) \cdots \mathbf{D}^{\beta_{m_B}} u(t) \\ -\mathbf{D}^{\alpha_1} y^*(t) \cdots -\mathbf{D}^{\alpha_{m_A}} y^*(t) \end{bmatrix} \quad (12)$$

and

$$\theta^T = [b_0 \ b_1 \ \dots \ b_{m_B} \ a_1 \ \dots \ a_{m_A}]. \quad (13)$$

Minimizing the  $L_2$  norm of  $\varepsilon(t)$ :

$$J = \int_0^T (\varepsilon(t))^2 dt, \quad (14)$$

with respect to  $\theta$ , leads to the following least squares estimates:

$$\hat{\theta}_{LS} = \left[ \int_0^T \varphi^*(t)^T \varphi^*(t) dt \right]^{-1} \int_0^T \varphi^*(t)^T y^*(t) dt. \quad (15)$$

Or, after a numerical discretization, by defining  $\mathbf{Y}$  as a column vector of the system output and  $\Phi$  as a regression matrix where the columns are fractional derivatives of the input and the output signals:

$$\Phi^* = [\varphi^*(T_s) \ \varphi^*(2T_s) \ \dots \ \varphi^*(KT_s)]^T, \quad (16)$$

$\hat{\theta}_{LS}$  can be approximated by:

$$\hat{\theta}_{LS} = (\Phi^{*T} \Phi^*)^{-1} \Phi^{*T} \mathbf{Y}^*. \quad (17)$$

Direct fractional differentiations of noisy output lead however to inaccurate results. Consequently, state variable filters are preferred to the direct least squares estimates.

### 2.2 Least Squares-based SVF method (lssvf)

As in the integer case, fractional differentiation of noisy signals amplifies the noise. Hence, a linear transformation (low-pass filter) is applied to input and output signals (Cois et al. (2001)). Poisson's filters can, for example, be chosen such as:

$$F_\gamma(s) = \frac{s^\gamma}{\left(\left(\frac{s}{\omega_c}\right)^\gamma + 1\right)^N}. \quad (18)$$

The order  $N$  is usually chosen such that  $\gamma N > \alpha_{m_A}$ . The basic idea in using Poisson's filter is to obtain the behavior of a differentiator in low frequencies and to filter noise in high frequencies as shown in the Bode diagram of Fig. 2. However, this filter introduces a phase shift near the cut-off frequency. The approximation is usually good enough to initialize the *srivcf* algorithm. Hence, differentiated filtered input  $\mathbf{D}^{\beta_i} u_f$  and output  $\mathbf{D}^{\alpha_j} y_f$  signals are obtained at the output of the filters (18):

$$\begin{cases} \mathbf{D}^{\beta_i} u_f(t) = u(t) * \mathcal{L}^{-1}\{F_{\beta_i}(s)\} \\ \mathbf{D}^{\alpha_j} y_f(t) = y^*(t) * \mathcal{L}^{-1}\{F_{\alpha_j}(s)\}, \end{cases} \quad (19)$$

where  $*$  and  $\mathcal{L}^{-1}$  stand respectively for the convolution operator and the inverse Laplace transform.

Instead of minimizing the  $L_2$  norm of  $\varepsilon(t)$  as in (11), the  $L_2$  norm of  $\varepsilon_f(t)$  is now minimized:

$$\varepsilon_f(t) = y_f^*(t) - \varphi_f^*(t)^T \theta, \quad (20)$$

where

$$\varphi_f^*(t)^T = \begin{bmatrix} \mathbf{D}^{\beta_0} u_f(t) \cdots \mathbf{D}^{\beta_{m_B}} u_f(t) \\ -\mathbf{D}^{\alpha_1} y_f^*(t) \cdots -\mathbf{D}^{\alpha_{m_A}} y_f^*(t) \end{bmatrix}. \quad (21)$$

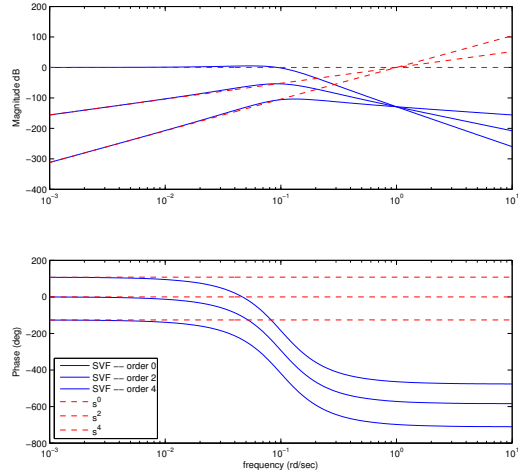


Fig. 2. Bode diagram of the fractional state variable filters

The solution is approximated (due to the integral) by the classical least squares-based SVF as in (17):

$$\hat{\theta}_{LSSVF} = (\Phi_f^{*T} \Phi_f^*)^{-1} \Phi_f^{*T} \mathbf{Y}_f^*, \quad (22)$$

where

$$\Phi_f^* = [\varphi_f^*(T_s) \ \varphi_f^*(2T_s) \ \dots \ \varphi_f^*(KT_s)]^T. \quad (23)$$

As in the integer case, Cois et al. (2001) showed that the least squares estimator (22) is biased in presence of noisy output due to the correlation between  $\varphi_f^*(t)$  and  $y^*(t)$ .

### 2.3 Instrumental Variable-based SVF method (ivsvf)

The estimator of the instrumental variable is a traditional variant of the least squares method. It presents the advantage of relying on linear regression techniques. The principle of the instrumental variable method relies on the so called instruments generally obtained from a parallel model  $y_f^{IV}$ . Hence, an instrumental variable regression vector is built:

$$\varphi_f^{IV}(t)^T = \begin{bmatrix} \mathbf{D}^{\beta_0} u_f(t) \cdots \mathbf{D}^{\beta_{m_B}} u_f(t) \\ -\mathbf{D}^{\alpha_1} y_f^{IV}(t) \cdots -\mathbf{D}^{\alpha_{m_A}} y_f^{IV}(t) \end{bmatrix}. \quad (24)$$

The conditions that the instrumental variable vector should satisfy, are the following (see e.g. Ljung (1999)) ( $\mathbf{E}[\cdot]$  stands for the mathematical expectation):

$$\begin{cases} \mathbf{E} [\varphi_f^{IV}(t) \varphi_f^{*T}(t)] \text{ is non singular} \\ \mathbf{E} [\varphi_f^{IV}(t) p(t)] = 0. \end{cases} \quad (25)$$

Hence, the instruments, gathered in  $\varphi_f^{IV}$ , need to be sufficiently correlated to all components of the regression vector  $\varphi_f^*$ , and totally uncorrelated to the output noise  $p(t)$ .

The IV-based solution is now approximated by:

$$\hat{\theta}_{IVSVF} = (\Phi_f^{IVT} \Phi_f^*)^{-1} \Phi_f^{IVT} \mathbf{Y}_f^*, \quad (26)$$

where

$$\Phi_f^{IV} = [\varphi_f^{IV}(T_s) \ \varphi_f^{IV}(2T_s) \ \dots \ \varphi_f^{IV}(KT_s)]^T. \quad (27)$$

2.4 Simplified refined instrumental variable for continuous-time fractional models (*srivcf*)

As shown by Young (1981), the optimal estimator (asymptotically without bias and of minimal variance) is obtained when the SVF filters  $F_\gamma(s)$  in (18) are replaced by:

$$F_\gamma^{opt}(s) = \frac{s^\gamma}{H(s)A(s)}, \quad (28)$$

where  $H(s)$  is the *real* noise model,  $A(s)$  the denominator of the transfer function of the real system; and the hypothetical noise free deterministic output  $y(t)$  is used when evaluating the derivatives of the output in (19). As a result, the optimal regression vector is obtained:

$$\varphi_f^{opt}(t)^T = \begin{bmatrix} \mathbf{D}^{\beta_0} u_f(t) \cdots \mathbf{D}^{\beta_{m_B}} u_f(t) \\ -\mathbf{D}^{\alpha_1} y_f(t) \cdots -\mathbf{D}^{\alpha_{m_A}} y_f(t) \end{bmatrix}. \quad (29)$$

The optimal estimator necessitates to know many characteristics of the system to be identified. It will be obtained only if the real system and noise models are known; however the optimality is not achievable in practical cases.

A stochastic approach proposed by Young (1981) and Young (2002) is the **Simplified Refined Instrumental Variable for Continuous systems** (*srivc*). This approach includes an adaptive pre-filtering based on the optimal solution when the noise  $p$  is a white gaussian one ( $H(s) = 1$ ). This method is a logical extension of the instrumental variable estimator with SVF. This estimator is extended to fractional models and is named *srivcf* (*srivc* for fractional models).

To optimize the instruments, it is required to implement an iterative algorithm. This algorithm can be initialized by the  $\hat{\theta}_{LSSVF}$  or  $\hat{\theta}_{IVSVF}$  estimates (22). Once the first estimation of the denominator is available, the new derivatives of input and output signals are computed by replacing  $F_\gamma(s)$  in (18) by:

$$F_\gamma^{(iter)}(s, \hat{\theta}) = \frac{s^\gamma}{\hat{A}(s)} = \frac{s^\gamma}{1 + \sum_{j=1}^{m_A} \hat{a}_j s^{\alpha_j}}, \quad (30)$$

with  $\hat{a}_j$  the estimated  $a_j$  coefficient and  $(iter) = 1, 2, \dots$  stands for the iteration number.

The instruments, the derivatives of the input and of the output, are computed at each iteration by:

$$\begin{cases} \mathbf{D}^{\beta_i} u_f(t) = u(t) * \mathcal{L}^{-1}\{F_{\beta_i}^{(iter)}(s)\} \\ \mathbf{D}^{\alpha_j} y_f^*(t) = y^*(t) * \mathcal{L}^{-1}\{F_{\alpha_j}^{(iter)}(s)\} \\ \mathbf{D}^{\alpha_j} y_f^{IV}(t) = y^{IV}(t) * \mathcal{L}^{-1}\{F_{\alpha_j}^{(iter)}(s)\}. \end{cases} \quad (31)$$

The regression vectors  $\varphi_f^*(t)$  and  $\varphi_f^{IV}(t)$  are formed as in (21) and (24). Finally, the parameter vector  $\hat{\theta}_{srivcf}^{(iter)}$  is computed at each iteration as:

$$\hat{\theta}_{srivcf}^{(iter)} = \left( \Phi_f^{IVT} \Phi_f^* \right)^{-1} \Phi_f^{IVT} \mathbf{Y}_f^*. \quad (32)$$

As in the rational case, the *srivcf* estimator is asymptotically unbiased whatever additive zero-mean noise is present. When the additive noise is white, *srivcf* has minimum variance and the asymptotic covariance matrix of the estimation errors associated with the estimate of  $\hat{\theta}_{srivcf}$  is given by:

$$\mathbf{P}_{\hat{\theta}} = \hat{\sigma}^2 (\Phi_f^{IVT} \Phi_f^{IV})^{-1}, \quad (33)$$

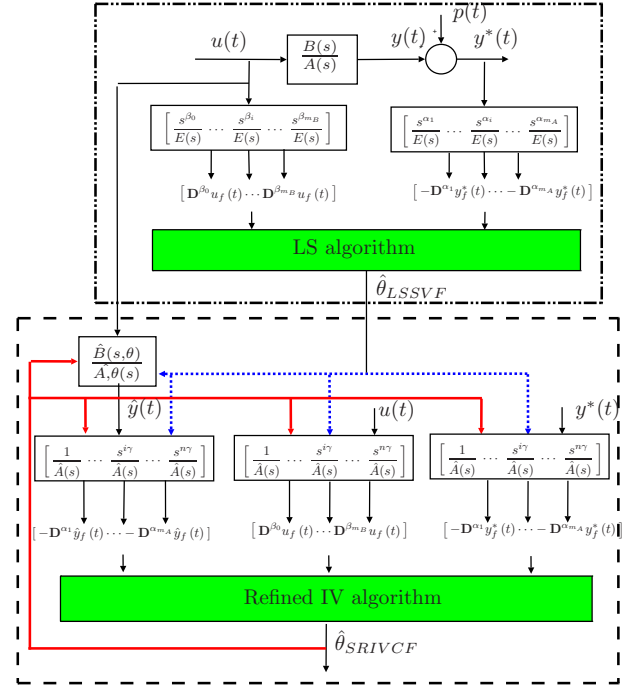


Fig. 3. Iterative optimal *srivcf*

where  $\hat{\sigma}^2$  is the empirical estimation of the noise variance.

2.5 Summary of the *srivcf* algorithm

The overall *srivcf* algorithm is summarized in this section and sketched in Fig. 3.

**Step 1** use the fractional state variable filter estimates (22) to generate an initial ( $iter = 1$ ) estimate of the transfer function model parameter vector  $\theta^{(1)}$ .

**Step 2** iterative instrumental variable estimation with prefilters.

**for**  $iter = 2$  to convergence

(i) Generate the instrumental variables  $y^{IV}$  from the auxiliary model with the estimated polynomials based on the estimated parameter vector  $\theta^{(iter-1)}$ .

(ii) Update the filter  $F_\gamma^{(iter)}(s, \hat{\theta})$  in (30) with the new estimated parameters. Then, evaluate the prefiltered derivatives of  $u(t)$ ,  $y(t)$  and  $y^{IV}(t)$  as in (31).

(iii) Based on these prefiltered data, compute the new estimates  $\theta^{(iter)}$  as in (32).

**end for**

**Step 3** compute the estimated parametric error covariance matrix associated with the parameter estimates with (33).

It may happen, when the bias is important, that the estimated parameters from Step 1 lead to unstable models. In this case, an empirical method can be used to stabilize unstable poles.

3. NUMERICAL EXAMPLE

3.1 System specifications

To illustrate the efficiency of this algorithm in system identification using rational and/or fractional models, the following system is simulated:

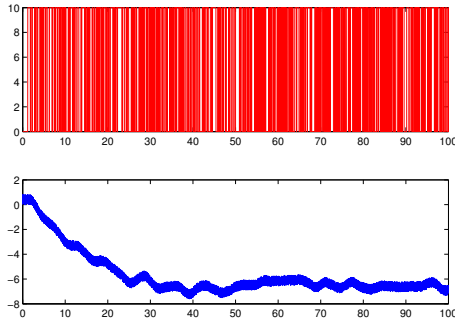


Fig. 4. Input and noisy output signals used for identification.

$$G_0(s) = \frac{K(-Ts^\nu + 1)}{\left(\left(\frac{s}{\omega_1}\right)^{2\nu} + 2\zeta_1\left(\frac{s}{\omega_1}\right)^\nu + 1\right)\left(\left(\frac{s}{\omega_2}\right)^{2\nu} + 2\zeta_2\left(\frac{s}{\omega_2}\right)^\nu + 1\right)}, \quad (34)$$

with  $\nu = 0.5$ ,  $K = -1$ ,  $T = 0.5$ ,  $\omega_1 = 0.2$  rad/s,  $\zeta_1 = -0.4$ ,  $\omega_2 = 1$  rad/s and  $\zeta_2 = -0.65$ .

This fractional system has two resonant modes: one at  $\omega_1 = 0.2$  rad/s with  $\zeta_1 = -0.4$  and another one at  $\omega_2 = 1$  rad/s with  $\zeta_2 = -0.65$  (for stable fractional systems, a resonant peak may occur for negative pseudo-damping factors  $\zeta_{1,2}$ <sup>2</sup>). Furthermore, the system has a positive  $s^\nu$ -zero at  $s^\nu = 2$ .

The input signal  $u(t)$ , a pseudo random binary sequence (prbs), is applied to the system (34). The prbs amplitude is intentionally chosen between 0 to 10 in order to be able to capture the rapid dynamics and the long memory behaviour (slow dynamics) typically present in fractional differentiation systems. The output,  $y(t)$ , is corrupted by an additive Gaussian white noise  $p(t)$  with zero mean and a signal to noise ratio (S/N) of 20dB as shown in Fig. 4. The sampling period is fixed to  $T_s = 10^{-2}$ s.

The instrumental variables are computed according to the iterative algorithm described in §2.5 in order to rippen the parametric estimation. The model, which parameters are to be estimated, is set to:

$$G(s) = \frac{b_1 s^\nu + b_0}{a_4 s^{4\nu} + a_3 s^{3\nu} + a_1 s^{2\nu} + a_1 s^\nu + 1}. \quad (35)$$

To be in the same class as the true system, the commensurable order must be set to 0.5. However, the commensurable order is not always known; when unknown, the optimal commensurable order can be estimated.

### 3.2 The choice of commensurable order

At the first stage, *srivcf* is applied with commensurable orders varying from  $\nu = 0.1$  to  $\nu = 1.9$  with a step of 0.05. The log-normalized cost function

$$J_{dB} = 10 \log_{10} \left( \frac{\sum_{k=1}^K (y_m(kT_s) - y^*(kT_s))^2}{\sum_{k=1}^K y_m(kT_s)^2} \right), \quad (36)$$

where  $y_m$  stands for the model output, is plotted versus the commensurable order in the region of interest, i.e.  $\nu \in [0.3, 1]$ , in Fig. 5. As expected, the best commensurable

<sup>2</sup> When  $\nu \neq 1$ ,  $\zeta$  cannot be considered as a damping factor.

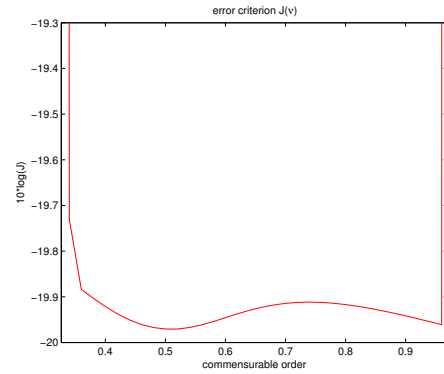


Fig. 5. Cost function in dB vs the commensurable order.

	true	srivcf ( $\nu = 0.5$ )		ivsvf	
		$E(\hat{\theta})$	$\sigma(\hat{\theta})$	$E(\hat{\theta})$	$\sigma(\hat{\theta})$
$a_4$	25	25.037	0.026	25.149	0.165
$a_3$	-36.5	-36.546	0.029	-36.671	0.106
$a_2$	31.2	31.235	0.015	31.252	0.034
$a_1$	-5.3	-5.302	0.004	-5.311	0.006
$b_1$	0.5	0.500	0.002	0.502	0.015
$b_0$	-1	-1.002	0.001	-0.999	0.002

Table 1. Comparison between *srivcf* and *ivsvf* methods,  $E(\hat{\theta})$  being the mean and  $\sigma(\hat{\theta})$  the standard deviation

order is at  $\nu = 0.5$ , accordingly to the true system (34).

### 3.3 Comparison between ivsvf and srivcf methods

This comparison will be done when the model (35) is in the same class as the true system, with  $\nu = 0.5$ . The objective here is to study the performance of the *srivcf* method as compared to the *ivsvf* method (Cois et al. (2001)). With different noise realizations with a S/N = 20dB, a Monte Carlo simulation of 200 runs has been performed. For each realization of noise, two models were evaluated: one with *ivsvf*, and another with *srivcf*. The Monte Carlo simulation results are displayed in Table 1.

Even though *ivsvf* and *srivcf* methods are both asymptotically unbiased (Young (2002)), *srivcf* significantly improves the parameter variance, as displayed in the  $\sigma(\hat{\theta})$  columns (standard deviation) of Table 1.

The frequency responses of the whole set of models obtained with *srivcf* and *ivsvf* methods are plotted in Fig. (6-a) and (6-b) respectively. As expected the 200 models obtained with *srivcf* fit exactly the Bode diagram of the simulated system (34) as compared to the *ivsvf* ones which have a larger dispersion.

### 3.4 Comparison between a fractional model and a rational one both evaluated with srivcf method

Considering *srivcf*, rational models are compared to fractional ones using a Monte Carlo simulation of 200 runs with S/N = 20dB. For each realization of noise, two models are evaluated: one for the best commensurable order  $\nu = 0.5$ , and the other for a rational model with  $\nu = 1$ .

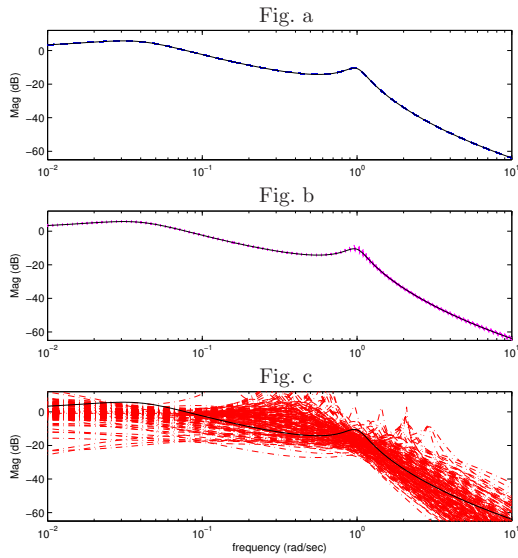


Fig. 6. Magnitude Bode diagrams of the models estimated for the 200 Monte Carlo simulation runs – Figure a: the true system (—), the fractional models ( $\nu = 0.5$ ) obtained with *srivcf* (--- lines) – Figure b: the true system (—), the fractional models ( $\nu = 0.5$ ) obtained with *ivsvf* (--- lines) – Figure c: the true system (—), and the rational models ( $\nu = 1$ ) obtained with *srivcf* (-.- lines).

For  $\nu = 0.5$ , the model coefficients  $b_i$  and  $a_j$  in (35) converged to the exact values (34) for all 200 Monte Carlo runs. However, for  $\nu = 1$ , unstable and inappropriate models are obtained. The frequency responses of the whole set of 200 rational models obtained for  $\nu = 1$  are plotted in Fig. (6-c). As expected, the rational models give less satisfactory results as the asymptotic slopes of rational systems are multiples of 20dB per decade in the magnitude Bode diagram, compared to fractional models which can have any asymptotic slope (Oustaloup (1995)). If a rational model equivalent to a fractional model in a desired frequency range were to be obtained, then a much higher order model would be required.

#### 4. CONCLUSION AND OUTLOOKS

In this paper, the simplified refined instrumental variable for continuous-time systems has been extended to fractional models (*srivcf*). The *srivcf* iterative algorithm is initialized from the traditional least squares-based SVF estimates. Numerical simulations have illustrated the performance of the proposed method.

The proposed approach can be seen to have a single tuning parameter: the commensurable order. As shown in the example, the cost function is tightly linked to the commensurable order. More studies are yet to be done on fractional models. Here, only the coefficients were estimated; the parametric estimation could be extended to the commensurable order. An interesting perspective would be to evaluate the procedure based on the properties of the instrumental product matrix (see Young and Jakeman (1980) or Young et al. (2006)) for identifying the commensurable model prior to parameter estimation. Another

interesting perspective is to develop a *rivcf* algorithm for hybrid Box-Jenkins models as in rational systems (Young et al. (2006)) in order to obtain minimum variance parameters in presence of colored noise.

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