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**A New Minimum Distance Estimation  
Procedure of ARFIMA Processes**

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# A New Minimum Distance Estimation Procedure of ARFIMA Processes

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## Abstract

A new parametric minimum distance time-domain estimator for ARFIMA processes is introduced in this paper. The proposed estimator minimizes the sum of squared correlations of residuals obtained after filtering a series through ARFIMA parameters. The estimator is easy to compute and is consistent, asymptotically normally distributed and efficient for fractionally integrated (FI) processes with an integration order  $d$  strictly greater than  $-0.75$ . Therefore, it can be applied to both stationary and non-stationary processes. Deterministic components are also allowed in the DGP. Furthermore, as a by-product, the estimation procedure provides an immediate check on the adequacy of the specified model. This is so because the criterion function, when evaluated at the estimated values, coincides with the Box-Pierce goodness of fit statistic. Empirical applications and Monte-Carlo simulations supporting the analytical results and showing the good performance of the estimator in finite samples are also provided.

*Keywords:* fractional integration; nonstationary long-memory time series; minimum distance estimation;

*JEL Classification:* C13, C22.

## 1. INTRODUCTION

A new estimation procedure for Autoregressive Fractionally Integrated Moving Average (ARFIMA) processes is proposed in this paper. First introduced by Granger and Joyeux (1980) and Hosking (1981), these processes have become very popular due to their ability in

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providing a good characterization of the long-run properties of many economic and financial time series. They are also very useful for modeling multivariate time series, since they can capture a larger number of long term equilibrium relations among economic variables than the traditional multivariate ARIMA models. See Baillie (1996) and Henry and Zaffaroni (2002) for a survey on this topic.

The estimator introduced in this paper belongs to the Minimum Distance (MD) class. The idea of the estimation procedure is quite simple: the parameters of the ARFIMA model are estimated by minimizing the sum of the squared autocorrelations of the residuals, obtained after filtering the original series through ARFIMA parameters. The proposed estimator is closely related to the MD estimator considered in Tieslau, Schmidt and Baillie (1996), and to the Adjusted MD estimator proposed in Chung and Schmidt (1995). Nevertheless, as it will be seen shortly, it presents important advantages over those estimators. It is denoted “Generalized Minimum Distance” (GMD) estimator since it extends previous approaches in this area to more general setups. In particular, the proposed estimator is easy to compute, has very good asymptotic and finite sample properties and is able to circumvent most of the problems present in the above-mentioned techniques. It can be applied to  $FI(d)$  series for values of  $d > -0.75$ , thus covering stationary as well as non-stationary ranges of  $d$ . Furthermore, it turns out to be asymptotically equivalent to ML estimation, (i.e.,  $\sqrt{T}$ -consistent, asymptotically normally distributed and efficient), for all  $d > -0.75$ .

This technique has been developed in the time domain, usually preferred in applied work. Relative to other time domain approaches, such as Maximum Likelihood estimation (MLE), it presents the additional advantage that it is not necessary to specify a particular distribution for the innovation process. Also, it is computationally faster than exact MLE for more complex ARFIMA( $p, d, q$ ) processes, since the evaluation of the likelihood requires finding the inverse of a  $T \times T$  matrix of population autocovariances, whose elements are complicated (hypergeometric) functions of the unknown parameters. For large sample sizes, accurate computation of such matrices could be a non-trivial task, even with today’s computing technology. As it will be seen, the computation of this estimator does not present these difficulties.

Frequency domain estimators are also very popular in this literature, mainly due to their computational simplicity and their good asymptotic properties, as it is the case of the Whittle estimator (see Fox and Taqqu, (1986) and Dahlhaus, (1989)). Nevertheless, the latter technique has also an important drawback with respect to the GMD estimator presented in this paper. The former estimator is consistent, asymptotically normal and efficient when the memory parameter  $d$  is known to lie in the stationary and invertible range and this restriction is correctly imposed. For the general case where  $d$  is completely unknown and possibly non-stationary, Velasco and Robinson (2000) showed that it is needed

to resort to tapered data to achieve consistency. Tapering increases the variance of the estimators and therefore induces an efficiency loss. This problem does not apply to the GMD estimator, since it is efficient for all  $d > -0.75$ . This implies that, in the general case where no information about  $d$  is available before estimation, GMD is more efficient than Whittle estimation.

Another interesting feature is that the estimation procedure proposed in this paper provides, as a by-product, an immediate check on adequacy of the specified parametric model. This is so because the criterion function, evaluated at the estimated values, yields the Box-Pierce goodness-of-fit statistic which has been largely used for these purposes in the literature.

It is also remarkable that the proposed framework can be easily extended to more general settings. For instance, the estimator can be easily robustified against conditional heteroscedasticity, just by introducing a modification in the definition of the residuals used to compute the autocorrelations. Also in this case it remains asymptotically equivalent to MLE. Finally, it can be extended, along the lines of Wright (1999), to the estimation of the fractionally integrated stochastic volatility model.

The rest of the paper is structured as follows. The ARFIMA model and the definition of the residuals are introduced in Section 2. The GMD estimation procedure and the asymptotic properties of the estimator are discussed in Section 3. The results of some simulation experiments, designed to evaluate the performance in finite samples of the proposed estimator, are described in Section 4. Section 5 derives the asymptotic distribution of residual autocorrelations and applies the result to the Box-Pierce (1970) and Box-Ljung (1978) goodness-of-fit statistics. Section 6 investigates their finite sample performance. An application of the described methods to empirical data is provided in Section 7. The conclusions of the paper are presented in Section 8. All proofs are gathered in Appendix 1. Appendix 2 incorporates the results of other Monte Carlo experiments.

The following conventional notation is adopted throughout the paper.  $L$  is the lag operator;  $\Delta = (1 - L)$ ;  $\Gamma(\cdot)$  denotes the gamma function; “ $\xrightarrow{w}$ ” and “ $\xrightarrow{p}$ ” denote weak convergence and convergence in probability, respectively;  $\|\cdot\|$  is the euclidean norm for  $m \times n$  matrices defined as  $\|A\| = \left(\sum_{i=1}^m \sum_{j=1}^n a_{ij}^2\right)^{1/2}$ ;  $\{\pi_i(d)\}_{i=0}^{\infty}$  represents the sequence of coefficients associated with the expansion of  $\Delta^d$  in powers of  $L$ , such that  $\Delta^d = \pi_0(d) + \pi_1(d)L + \pi_2(d)L^2 + \dots$ , and

$$\pi_i(d) = \frac{\Gamma(i-d)}{\Gamma(-d)\Gamma(i+1)}. \quad (1)$$

## 2. THE MODEL

The definition of ARFIMA process adopted in this article is similar to that used in Beran (1995) or Sowell (1990). Consider the ARFIMA( $p, d_0, q$ ) process  $y_t, t = 1, 2, \dots, T$ , that can be written as

$$\Phi_0(L) \Delta^{\varphi_0} (\Delta^{m_0} y_t - \mu_0) = \Theta_0(L) \varepsilon_t. \quad (2)$$

In the previous definition  $\{\varepsilon_t\}_{t=-\infty}^{\infty}$  is a sequence of i.i.d zero-mean random variables with unknown variance  $\sigma^2$  and finite fourth moment,  $E(\varepsilon_t^4) = \mu_4 < \infty$ .  $\Phi_0(L)$  and  $\Theta_0(L)$  are autoregressive and moving average polynomials of order  $p$  and  $q$ , respectively, with all roots outside the unit circle. Throughout, it will be assumed that  $p$  and  $q$  are known natural numbers. The memory parameter,  $d_0$ , belongs to the closed interval  $[\nabla_1, \nabla_2]$ , with  $-0.75 < \nabla_1 < \nabla_2 < \infty$ . It is composed as the sum of an integer and a fractional part such that  $d_0 = m_0 + \varphi_0$ . The integer  $m_0 = \lfloor d_0 + 1/2 \rfloor$ , where  $\lfloor \cdot \rfloor$  denotes integer part, is the number of times that  $y_t$  must be differenced to achieve stationarity (therefore  $m_0 \geq 0$ ). The parameter  $\varphi_0$ , the fractional part, lies in the interval  $(-0.75, 0.5)$ , in such a way that, for a given  $d_0$ ,  $\varphi_0 = d_0 - \lfloor d_0 + 1/2 \rfloor$ . Consequently, once the process  $y_t$  is differenced  $m_0$  times, the differenced process is a stationary fractionally integrated process with integration order  $\varphi_0$ . More explicitly, if  $x_t(m_0) = (\Delta^{m_0} y_t - \mu_0)$ , then  $x_t(m_0)$  is a (stationary)  $FI(\varphi_0)$ . For  $m_0 = 0$ ,  $\mu_0$  is the expected value of the stationary process  $y_t$  and for  $m_0 \geq 1$ ,  $\mu_0 \neq 0$  implies a deterministic polynomial trend. This definition is also similar, for instance, to that used in Velasco and Robinson (2000). For the asymptotic implications of using different definitions of ARFIMA processes see Marinucci and Robinson (1999).

To derive the new estimator we need to define the residuals of the process. For that purpose, we adopt Beran's (1995) definition of residuals and provide two alternative expressions according to whether the mean,  $\mu_0$ , is known or unknown.

### 2.1. Residuals when $\mu_0$ is known and equal to zero.

Let  $\psi = (\phi_1, \dots, \phi_p, \theta_1, \dots, \theta_q)' \in \mathfrak{R}^{p+q}$  be the vector containing the autoregressive and moving average parameters and  $\lambda = (d, \psi) \in \mathfrak{R}^{p+q+1}$ . Also let  $\lambda_0 = (d_0, \psi_0)'$  represent the vector containing the true parameter values. Analogously, define  $\lambda^* = (\varphi, \psi)'$  and  $\lambda_0^* = (\varphi_0, \psi_0)'$ . The (infinite)<sup>1</sup> autoregressive representation of  $y_t$  is given by

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<sup>1</sup>Notice that this expansion is valid for all  $d_0 > -1$ . For values of  $d_0 > -0.5$ , this is a well known result due to Hosking (1981). When  $d_0 \in (-1, -0.5)$ , Odaki (1993, Theorem 2) shows that although the coefficients  $\pi_j(d_0)$  are not square summable and, consequently, the same applies to the coefficients  $\alpha_j(\lambda_0)$ , the process is still invertible and therefore the autoregressive inversion is well defined.

$$\sum_{j=0}^{\infty} \alpha_j(\lambda_0) y_{t-j} = \varepsilon_t, \quad (3)$$

where  $\{\alpha_j(\lambda_0)\}_{j=0}^{\infty}$  are the coefficients associated to the expansion of  $\Phi_0(L)\Theta_0(L)^{-1}\Delta^{d_0}$  in powers of  $L$ . Given the observations  $y_1, \dots, y_T$ , the innovations  $\varepsilon_t$  cannot be computed directly, since an infinite sample would be needed. Nevertheless, they may be estimated by

$$e_t(\lambda) = \sum_{j=0}^{t-m-1} \alpha_j(\lambda) y_{t-j}, \quad t = 1, \dots, T. \quad (4)$$

## 2.2. Residuals when $\mu_0$ is unknown.

When  $\mu_0$  is unknown the residuals defined above need to be adjusted. Again, following Beran (1995), we consider:

$$x_t(m) = \Delta^m y_t$$

and

$$\bar{x}(m) = \frac{1}{T-m} \sum_{t=m+1}^T x_t. \quad (5)$$

Since  $x_t(m_0)$  is stationary and ergodic, the sample mean  $\bar{x}(m_0)$  is a consistent estimator of  $\mu_0$ . Therefore, adjusted residuals can be defined as:

$$e_t(\lambda) = \sum_{j=0}^{t-m-1} \alpha_j(\lambda^*) (x_{t-j}(m) - \bar{x}(m)), \quad t = m+1, \dots, T. \quad (6)$$

where  $\{\alpha_j(\lambda^*)\}_{j=0}^{\infty}$  are the coefficients associated to the expansion of  $\Phi(L)\Theta(L)^{-1}\Delta^\varphi$  in powers of  $L$ .

## 3. GENERALIZED MINIMUM DISTANCE ESTIMATION OF ARFIMA PROCESSES.

Minimum Distance (MD) is a classical estimation approach in the econometric literature. This technique encompasses other very popular procedures such as Generalized Method of Moments (GMM), Non-Linear Least Squares (NLS) or Maximum Likelihood (ML) among others. In a general framework, this technique would work as follows: If  $\lambda_0 \in \Lambda$  is the vector of parameters of interest, where  $\Lambda$  is the set of possible parameter values, and  $y_t$  is the available data, MD estimation provides a class of estimators that minimize the following criterion function,

$$V_T(\lambda) = \hat{g}_T(\lambda)' \hat{W} \hat{g}_T(\lambda), \quad (7)$$

where  $\hat{g}_T(\lambda)$  is a function of the data,  $y_t$ , and the parameters of interest,  $\lambda$ , that has to verify  $\hat{g}_T(\lambda_0) \xrightarrow{p} 0$ ;  $\hat{W}$  is a positive definite weighting matrix that defines the distance.

Under the standard regularity conditions, it can be proved that the resulting estimators are  $\sqrt{T}$ -consistent and asymptotically normally distributed (see, for instance, Newey and McFadden, 1994). Different choices of the function  $\hat{g}_T(\lambda)$  will generate different estimators. For instance, if  $\hat{g}_T(\lambda) = T^{-1} \sum_{t=1}^T g(y_t, \lambda)$ , where  $E(g(y_t, \lambda_0)) = 0$ , the minimization of the criterion function in (7) would provide a GMM estimator. The function  $\hat{g}_T(\lambda)$  considered in this paper can be interpreted as the difference between the sample and the population autocorrelations of the residuals defined in (4) or in (6). As it will be proved below, the function  $\hat{g}_T(\lambda)$  defined in this way fulfils the above mentioned requirements. With respect to the choice of  $\hat{W}$ , it is a well-known result that if  $\text{var}(\sqrt{T}\hat{g}_T(\lambda_0)) \xrightarrow{p} C$ , then the efficient weighting matrix,  $W_e$ , is given by  $W_e = C^{-1}$ , since in this case the asymptotic variance-covariance matrix of  $\hat{\lambda}$  simplifies to  $(J'_{\lambda_0} C^{-1} J_{\lambda_0})^{-1}$ , where  $J_{\lambda_0}$  is the limit of the Jacobian matrix of  $\hat{g}_T$  (see Newey and McFadden, 1994). Therefore, this would be in general a good choice. Some examples of  $\hat{W}$  corresponding to particular situations will be provided below. As it becomes clear from the discussion above, one of the main advantages of MD relative to ML estimation is that the former do not require to assume a particular distribution of the innovation sequence at any stage.

Several parametric and semiparametric MD techniques can be found in the literature of fractionally integrated (*FI*) processes. Robinson (1994a) proposed a semiparametric time domain procedure that exploits the fact that autocovariances in *FI* models are proportional to  $k^{2d-1}$  for large  $k$ . Robinson's estimator minimizes the expression:  $\sum_{k=n}^{p+n} (\hat{\gamma}_k - ck^{2d-1})^2$ . Hall et al. (1997) have analyzed the rates of convergence of this estimator but its distributional properties remain to be determined. Of particular relevance to this paper is the MD estimator proposed by Tieslau, Schmidt and Baillie (1996), henceforth TSB. They introduced a parametric time domain MD estimator that minimizes a distance between the estimated and the theoretical autocorrelations of an ARFIMA  $(p, d, q)$  process:

$$\hat{\lambda} = \arg \min_{\lambda \in \Lambda} (\hat{\rho}_{ky} - \rho_{ky}(\lambda))' \hat{W} (\hat{\rho}_{ky} - \rho_{ky}(\lambda)), \quad (8)$$

where  $\hat{\rho}_{ky}$  is the sample autocorrelation function of the (stationary) process  $y_t$  up to lag  $k$  (for a fixed value of  $k$ ),  $\rho_{ky}(\lambda)$  is the theoretical autocorrelation of the corresponding ARFIMA $(p, d, q)$  process up to the same lag and  $\hat{W}$  is a positive definite weighting matrix. The asymptotic optimal weighting matrix is  $W_e = C^{-1}$ , where  $C$  is the asymptotic variance-covariance matrix of  $\sqrt{T}\hat{\rho}_{ky}$ . Then, a suitable choice for  $\hat{W}$  in this context would be a consistent estimator of  $C^{-1}$ . Although theoretically very appealing, there remain significant problems with this procedure. First, it is restricted to stationary series since it requires the existence of autocorrelations. Second, it is  $\sqrt{T}$ -consistent and asymptotically normal only for  $d_0 < 0.25$ , due to the non-standard behavior of sample autocorrelations of ARFIMA processes outside this range. Moreover, the procedure is not efficient over its entire domain.

Chung and Schmidt (1995) have introduced a modification (Adjusted Minimum Distance Estimator) to the previous estimator. They have demonstrated, by applying the results on autocorrelations of Hosking (1996), that it is possible to obtain a  $\sqrt{T}$ -consistent and asymptotically normally distributed estimator of  $d$  in the whole invertible and stationary range,  $-0.5 < d < 0.5$ , if some functions of the autocorrelations are employed in the criterion function. Yet, it still remains not efficient and it is computationally almost as demanding as exact ML (see Sowell, 1992) since it requires the computation of the autocorrelations as functions of the unknown parameters. Along the same lines, Wright (1999) has proposed an estimator for the fractionally integrated stochastic volatility model and has proved that it is  $\sqrt{T}$ -consistent and asymptotically normally distributed only when  $d < 0.25$ . Galbraith and Zinde-Walsh (1997) have presented a parametric time-domain estimator based on an autoregressive approximation. This estimator can be applied to nonstationary series, since the existence of autocorrelations is not required, but its consistency has not been proved yet in this general framework. Other interesting references on the parametric estimation of possibly non stationary ARFIMA processes are Beran (1995), Tanaka (1999), Velasco and Robinson (2000), and Ling and Li (1997). See also Hauser (1999) for a finite sample comparison among ML estimation techniques.

Let us now describe the GMD estimator proposed in this paper. Consider the sample  $i$ th-autocorrelation associated with the residuals defined in (4) or in (6) given by:

$$\hat{\rho}_{e(\lambda)}(i) = \frac{\sum_{t=1}^{T-i} e_t(\lambda)e_{t+i}(\lambda)}{\sum_{t=1}^T e_t(\lambda)^2}. \quad (9)$$

Also define the vector  $\hat{\rho}_{ke(\lambda)}$  that contains the first  $k$  autocorrelations of the residuals :

$$\hat{\rho}_{ke(\lambda)} = \left( \hat{\rho}_{e(\lambda)}(1), \dots, \hat{\rho}_{e(\lambda)}(k) \right)'. \quad (10)$$

The following theorem will be very useful in the derivation of the subsequent theory:

**Theorem 1** *Consider the vector defined in (10) evaluated at  $\lambda = \lambda_0$ . Under the assumptions of Section 2, then:*

$$\left( \hat{\rho}_{ke(\lambda_0)} - \hat{\rho}_{k\varepsilon} \right) = o_p(1), \text{ for } d_0 > -1,$$

and

$$\sqrt{T} \left( \hat{\rho}_{ke(\lambda_0)} - \hat{\rho}_{k\varepsilon} \right) = o_p(1), \text{ for } d_0 > -0.75. \quad (11)$$

where  $\hat{\rho}_{k\varepsilon} = \left( \hat{\rho}_{\varepsilon}(1), \dots, \hat{\rho}_{\varepsilon}(k) \right)'$  is the vector that contains the sample first  $k$  autocorrelations (for fixed  $k$ ) associated to  $\varepsilon_t$ .

The previous theorem implies that the asymptotic distribution of the sample correlations of the residuals  $e_t(\lambda_0)$  associated to the ARFIMA process,  $y_t$ , coincides with that of the

sample correlations associated to the true innovations of  $y_t$ ,  $\varepsilon_t$ , for  $d_0 > -0.75$ . This means that the effect of the truncation is asymptotically negligible and therefore

$$\sqrt{T}\hat{\rho}_{ke(\lambda_0)} \xrightarrow{w} N(0, I_k), \text{ for } d > -0.75,$$

where  $I_k$  is the identity matrix of order  $k$ , with fixed  $k$ . Appendix 2 includes the results of some Monte Carlo experiments that investigate the small sample behavior of this approximation.

Following the argument in TSB (1996), we consider the minimization of a distance between the estimated and theoretical autocorrelations but, in place of the original series, the correlations of the above defined residuals are considered. Since the asymptotic variance of  $\sqrt{T}\hat{\rho}_{ke(\lambda_0)}$  is given by  $I_k$ , the identity matrix of order  $k$ , (Theorem 1), it follows (see (7)) that the efficient weighting matrix is  $I_k$ . Moreover, since  $\hat{\rho}_{ke(\lambda_0)} \xrightarrow{p} 0$ , the MD criterion function,  $V_{ke}$ , becomes

$$V_{ke}(\lambda, y) = \hat{\rho}'_{ke(\lambda)} \hat{\rho}_{ke(\lambda)} = \sum_{i=1}^k \hat{\rho}_{e(\lambda)}(i)^2, \quad (12)$$

and the GMD estimator  $\hat{\lambda}_k$  is defined as:

$$\hat{\lambda}_k = \arg \min_{\lambda \in \Lambda} V_{ke}(\lambda, y). \quad (13)$$

Notice that, since all the sample autocorrelations converge to zero and the efficient weighting matrix is the identity matrix, the criterion function simplifies notably with respect to that defined in (8), corresponding to the TSB approach. Also, it just requires the existence of the autocorrelations of the residuals but not those of the original series and therefore it can also be applied to non-stationary series, in contrast to the TSB estimator.

So far we have considered that the number of autocorrelations included in the criterion function,  $k$ , is a fixed number, as the above mentioned MD approaches have done. A further novelty of this paper consists of allowing  $k$  to grow to infinity with the sample size  $T$ . As it will be seen in Theorems 3 and 4, estimators computed with a fixed  $k$  are consistent and asymptotically normal, as long as  $k$  is greater than the  $p + q + 1^2$ . But, considering that  $k$  grows with  $T$  (Theorem 5) allows to obtain an asymptotically efficient estimator. The following theorem extends the results of Theorem 1 by showing that the sum of (weighted) correlations also converges to a Normal distribution, when the number of included correlations is allowed to go to infinity.

**Theorem 2** *Let  $\hat{\rho}_{ke(\lambda_0)}$  be the vector in (10), evaluated at  $\lambda = \lambda_0$ . Consider the series of  $h \times k$  stochastic matrices  $\{\hat{A}_k\}$  for  $k = 1, \dots, K$ , and let  $\{A_k\}_{k=1}^K$  be the probability limit of*

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<sup>2</sup>This restriction should be imposed to achieve identifiability of the parameters.

the former series. Now let  $k = k(T)$  be a function of  $T$  such that  $\lim_{T \rightarrow \infty} k(T) = \infty$  and  $\lim_{T \rightarrow \infty} k(T)/T = 0$ . If  $\|\hat{A}_k - A_k\| = o_p(k^{-1/2})$  and  $\lim_{k \rightarrow \infty} A_k A_k'$  is a bounded, positive definite matrix then,

$$\sqrt{T} \hat{A}_k \hat{\rho}_{ke(\lambda_0)} \xrightarrow{w} N(0, \quad).$$

The matrix  $\hat{A}_k$  contains the (stochastic) weights of the autocorrelation function. Later in the paper, this theorem will be used to show the asymptotic normality of the estimator when  $k$  tends to  $\infty$ . For that purpose,  $\hat{A}_k$  and  $A_k$  will be replaced by the Jacobian matrix of  $\hat{\rho}_{ke(\lambda)}$  and its corresponding probability limit, respectively.

As mentioned above, Beran (1995), Tanaka (1999) and Velasco and Robinson (2000) also present parametric estimators of both stationary and nonstationary fractionally integrated processes. The first two consider time-domain approximate Gaussian ML estimates, irrespective of the degree of nonstationarity. However, Beran's proof of consistency encounters some problems as Velasco and Robinson (2000) point out. The difficulties arise in the Taylor formula employed in the proof, since the  $o_p(1)$  term in the expansion of  $\sqrt{T}(\hat{\theta}_T - \theta_0)$  is only justified when  $\hat{\theta}_T$  lies in a sufficiently small neighborhood of  $\theta_0$ , which in turn presumes the consistency to be proved (see for instance, Brockwell and Davis, 1993, proposition 6.1.6). Nevertheless, the reported simulations support Beran's conclusions that this estimator is  $\sqrt{T}$ -consistent, asymptotic normally distributed and efficient for all  $d > -0.5$ . Tanaka (1999) extends Beran's approach to cover also non-invertible series. The estimates are found by minimizing the variance of the residuals but, since this variance is not a consistent estimator of the innovation variance for non-invertible FI series with  $d < -1$ , (see Odaki, 1993), Tanaka's estimator is not consistent either for  $d < -1$ . In the frequency domain, Velasco and Robinson (2000) show that the Whittle estimate is still appropriate for non-stationary processes as long as enough tapering is applied to the data. This induces an efficiency loss, since the introduction of tapering increases the variance of the estimators. There are other approaches in the frequency-domain literature that do not need to resort to tapering to achieve consistency in the non-stationary range of values of  $d$ , as is the case of the semiparametric modified log periodogram regression estimator (see Phillips, 1999).

The following theorems state the asymptotic properties of the GMD estimator proposed here. Theorems 3 and 4 state that it is consistent and asymptotically normally distributed, respectively, when  $T \rightarrow \infty$  and  $k$ , the number of correlations included in the criterion function, is a fixed number greater than  $p + q + 1$ . Theorem 5 extends the previous results to the case where  $k$  is a function of the sample size that also goes to infinity with  $T$ . It is shown that the estimator is not only consistent and asymptotically normally distributed in this case, but also that it is efficient since its asymptotic variance-covariance matrix is the inverse of the Fisher information matrix as it is of ML estimators, as established by Li and McLeod (1986) and Dahlhaus (1989). As in Beran (1995), Tanaka (1999) and Phillips

(2000), and in contrast to Velasco and Robinson (2000), no tapering is required to achieve these results.

**Theorem 3** *Let  $y_t$  be an ARFIMA( $p, d_0, q$ ) process under the hypotheses of Section 2. Also let  $\lambda_0$  be an interior point of the compact set  $\Lambda$ . Then, as  $T$  tends to infinity, it holds that*

$$\hat{\lambda}_k \xrightarrow{p} \lambda_0,$$

where  $\hat{\lambda}_k$  is the GMD estimator defined in (13) and  $k$  is a fixed number such that  $k \geq p + q + 1$ .

**Theorem 4** *Under the hypotheses of the previous theorem, it holds that:*

$$\sqrt{T}(\hat{\lambda}_k - \lambda_0) \xrightarrow{w} N(0, \Xi_k^{-1}), \quad (14)$$

(for definition of  $\Xi_k^{-1}$ , see Appendix 1).

The following theorem presents analogous results when the number of correlations to be included in the criterion function is not a fixed number but it grows with the sample size. By allowing for an increasing  $k$ , efficiency is achieved.

**Theorem 5** *Let  $y_t$  be an ARFIMA( $p, d_0, q$ ) process under the hypotheses of Section 2 and  $\lambda_0$  an interior point of the compact set  $\Lambda$ . Let  $\hat{\lambda}$  be the GMD estimator defined in (13) where the number of correlations included in the criterion function,  $k$ , is a function of  $T$  such that  $\lim_{T \rightarrow \infty} k(T) = \infty$  and  $\lim_{T \rightarrow \infty} k(T)/T = 0$ . Then, as  $T$  tends to infinity, it holds that*

$$\hat{\lambda} \xrightarrow{p} \lambda_0,$$

and

$$\sqrt{T}(\hat{\lambda} - \lambda_0) \xrightarrow{w} N(0, \Xi^{-1}), \quad (15)$$

where  $\Xi = \lim_{k \rightarrow \infty} \Xi_k$  is the Fisher information matrix (for definition, see Appendix 1).

The matrix  $\Xi$  is identical to the Fisher information matrix for ARMA processes except for the first row and column. Then, estimation of the remaining columns and rows is identical to that of ARMA processes (see, for instance, Brockell and Davis, 1993). See also Tanaka (1999) for a description of the estimation of the whole variance-covariance matrix.

The estimation approaches that just consider the invertible and stationary range of values of  $d$  ( $d \in (-0.5, 0.5)$ ) should first determine  $m_0$  in an exploratory way. In the case where the value of  $m_0$  is correctly guessed, an efficient estimator of this type and the estimator proposed in this paper will share the same asymptotic variance-covariance matrix and therefore they will be (asymptotically) equivalent (notice that the asymptotic covariance

matrix of  $\hat{\lambda}$  and  $\hat{\lambda}_k$  is independent on the value of  $d_0$ ). But when the assumption about the value of  $m_0$  is wrong, the former family of estimators will become inconsistent while this problem will not affect the GMD estimator. For this reason it is safer to use a method that covers the whole range of values of  $d$  and there is no loss of efficiency by doing so if the GMD estimator is employed. Nevertheless, when other methods are employed, such as the Whittle estimator with tapered data, some efficiency loss can occur.

On the other hand, the Box-Jenkins methodology shares the same exploratory approach for choosing  $m_0$  and, once this value has been chosen, inference on the rest of the ARMA parameters is carried out as if the differencing order was known a priori. This poses an additional problem since this procedure clearly underestimates the uncertainty of the problem and may lead to unrealistic confidence intervals for the remaining parameters.

#### 4. BEHAVIOR OF THE GMD ESTIMATOR IN FINITE SAMPLES.

In this section, a Monte Carlo study is conducted to investigate the small-sample performance of the MD estimator defined in (13). Processes of the form  $\Delta^{\varphi_0} (\Delta^{m_0} y_t - \mu_0) = u_t$  were generated, with four different specifications for  $u_t$ , namely  $u_t = \varepsilon_t$ ,  $u_t = \phi_1 u_{t-1} + \varepsilon_t$ ,  $u_t = (1 + \theta_1 L) \varepsilon_t$ , and  $u_t = \phi_1 u_{t-1} + \phi_2 u_{t-2} + \varepsilon_t$ ,  $\varepsilon_t \sim NID(0, 1)$  in all cases. Further,  $\mu_0$  is set equal to zero but the estimation procedure is carried out both considering that its value is known (and equal to zero) and also that it is unknown.

Before estimating the above-mentioned models, it is necessary to select the number of correlations ( $k$ ) to be included in the criterion function  $V_{ke}$ . Although increasing  $k$  always improves efficiency asymptotically, the suitable choice of  $k$  in finite-sample applications depends on the sample size  $T$  and both on the number ( $p + q + 1$ ) and the values of the parameters. Asymptotic theory does not help very much with respect to the right choice of  $k$  and, therefore, this is a question that should be addressed via Monte Carlo simulation. Table 1 presents the bias and the square root of the mean square error (SRMSE) of the GMD estimator in the case where  $u_t = \varepsilon_t$ , for two different sample sizes,  $T = 100$  and  $T = 400$ . The number of replications was 1000. Different values of  $k$  were used, namely, the closest integer to the quantities:  $T^{1/4}$ ,  $T^{1/3}$ ,  $T^{1/2}$  (more precisely  $k = 3, 5, 10$  and  $k = 4, 7, 20$  for  $T = 100, 400$  respectively). Table 2 presents analogous results for the case where  $\mu_0$  is unknown and has to be estimated. Tables 3, 4 and 5 display the figures obtained from similar experiments for the case where  $u_t$  is an  $AR(1)$ ,  $MA(1)$  or an  $AR(2)$  process with parameters  $\phi_1 = 0.6$ ,  $\theta_1 = 0.5$  or  $\phi_1 = 0.65$  and  $\phi_2 = -0.6$  respectively, and  $k = T^{1/4}$ . The value of the parameters have been chosen in order to facilitate comparison with the ones used in previous studies. Other values of  $k$  were also tried and they are not reported for economy of space but they are available upon request.

Figures in Tables 1 to 5 show the good performance of the proposed method. From Tables

1 and 2 it can be seen that the GMD estimator is surprisingly robust across different values of  $k$ . Moderate values of  $k$  provide the best results. For  $T = 100$ , the SRMSE is in general smaller for those estimates computed with  $k = T^{1/4}$ , although for  $T = 400$  the SRMSE is very similar and even smaller for those computed with  $T^{1/3}$ . With respect to efficiency, the asymptotic standard deviations for the ARFIMA(0,  $d$ , 0) case given in (14) are  $\pi^{-1}\sqrt{(6/T)}$ , which equals 0.078 and 0.039 for  $T = 100$  and 400, respectively. The reported SRMSE in Table 1 are a reasonable approximation to these values. It is also remarkable the small bias of the estimates, even for very large values of  $d$ . These values are usually negative which suggests that the GMD method slightly underestimates the memory parameter. Also note that the figures for the case where  $\mu_0$  is unknown do not differ significantly from the case where it is known. For the AR and MA cases different values of  $k$  were used and similar results were obtained. Again, the estimator performs better in terms of MSE when moderate values of  $k$  are used in the criterion function. The estimates in Table 3 to 5 have been calculated with  $k = T^{1/4}$ . The asymptotic variance-covariance matrix (*Asymp. Var*) for the ARFIMA(1,  $d$ , 0) process is given by

$$\text{Asymp. Var} \begin{pmatrix} \hat{d} \\ \hat{\phi}_1 \end{pmatrix} = \frac{1}{T} \begin{pmatrix} \frac{\pi^2}{6} & -\frac{1}{\phi_1} \log(1 - \phi_1) \\ -\frac{1}{\phi_1} \log(1 - \phi_1) & \frac{1}{1 - \phi_1^2} \end{pmatrix}^{-1}, \quad (16)$$

which delivers, for a value of  $\phi_1 = 0.6$ , asymptotic standard deviations equal to 0.256 and 0.262 for  $T = 100$ , and 0.128 and 0.131 for  $T=400$ , corresponding to  $\hat{d}$  and  $\hat{\phi}_1$  respectively. As for the ARFIMA(0,  $d$ , 1) case, asymptotic standard deviations can be computed from an analogous expression and, for  $\theta_1 = 0.5$ , they are equal to 0.222 and 0.246 for  $T = 100$  and 0.1108, and 0.1231 for  $T = 400$  corresponding also to  $\hat{d}$  and  $\hat{\theta}_1$  respectively. Again it is seen that the reported *SRMSEs* are a reasonable approximation to these values.

Table 6 compares the *SRMSE* of different estimators of  $d$  for the ARFIMA(0,  $d$ , 0) case. More specifically we consider the Whittle estimator (with Zhurbenko taper of order 2) proposed by Velasco and Robinson (VR), the ML estimators proposed by Sowell (SOW) and Beran (BER) and the minimum distance estimators by Tieslau et al. (TSB) and by Galbraith and Zinde-Walsh (GZW). The DGP used in this experiment was a fractional white noise with known mean equal to zero. The missing values in Table 6 stem from methods that are not defined for the whole range of values of  $d$ . It can be observed that the GMD estimator behaves similarly to the ML estimators and better than the remaining ones. It is also remarkable the good performance of the GMD estimator in the range of values of  $d$  in which other estimators are not defined.

**Table 1.** Estimation of  $d$  for the  $ARFIMA(0, d, 0)$  case

		$\mu_0$ known ( $\mu_0 = 0$ )								
$d_0$		-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$										
bias $\hat{d}$	$k = T^{1/4}$	-0.018	-0.032	-0.021	-0.017	-0.031	-0.019	-0.019	-0.026	-0.021
	$k = T^{1/3}$	-0.020	-0.031	-0.022	-0.029	-0.035	-0.021	-0.21	-0.028	-0.023
	$k = T^{1/2}$	-0.022	-0.032	-0.024	-0.032	-0.036	-0.023	-0.032	-0.030	-0.025
SRMSE $\hat{d}$	$k = T^{1/4}$	0.101	0.105	0.094	0.101	0.0970	0.093	0.093	0.104	0.098
	$k = T^{1/3}$	0.111	0.110	0.111	0.109	0.114	0.103	0.103	0.118	0.102
	$k = T^{1/2}$	0.116	0.117	0.114	0.113	0.122	0.116	0.118	0.122	0.121
$T = 400$										
bias $\hat{d}$	$k = T^{1/4}$	-0.003	-0.007	-0.008	-0.010	-0.008	0.008	0.005	0.006	0.008
	$k = T^{1/3}$	-0.003	-0.008	-0.010	-0.012	0.008	0.010	0.006	0.006	0.002
	$k = T^{1/2}$	-0.003	-0.008	-0.010	-0.013	0.010	0.010	0.007	0.008	0.009
SRMSE $\hat{d}$	$k = T^{1/4}$	0.045	0.042	0.045	0.041	0.042	0.046	0.044	0.044	0.046
	$k = T^{1/3}$	0.044	0.041	0.043	0.041	0.042	0.046	0.043	0.044	0.044
	$k = T^{1/2}$	0.047	0.042	0.045	0.042	0.044	0.048	0.044	0.045	0.046

**Table 2.** Estimation of  $d$  for the  $ARFIMA(0, d, 0)$  case

		$\mu_0$ unknown ( $\mu_0 = 0$ )								
$d_0$		-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$										
bias $\hat{d}$	$k = T^{1/4}$	-0.005	-0.030	0.002	-0.003	-0.026	-0.022	-0.024	-0.027	-0.020
	$k = T^{1/3}$	-0.007	-0.036	-0.002	-0.031	-0.027	-0.025	-0.026	-0.030	-0.022
	$k = T^{1/2}$	-0.009	-0.034	-0.027	-0.031	-0.030	-0.027	-0.029	-0.029	-0.025
SRMSE $\hat{d}$	$k = T^{1/4}$	0.101	0.107	0.097	0.103	0.104	0.100	0.102	0.104	0.099
	$k = T^{1/3}$	0.113	0.107	0.106	0.109	0.110	0.114	0.104	0.107	0.105
	$k = T^{1/2}$	0.117	0.109	0.111	0.112	0.123	0.117	0.117	0.117	0.109
$T = 400$										
bias $\hat{d}$	$k = T^{1/4}$	0.008	-0.007	0.001	-0.05	-0.008	-0.010	-0.004	-0.009	-0.008
	$k = T^{1/3}$	0.009	-0.009	0.003	-0.007	-0.007	-0.010	-0.009	-0.010	-0.010
	$k = T^{1/2}$	0.022	-0.009	-0.001	-0.007	-0.008	-0.011	-0.009	-0.012	-0.016
SRMSE $\hat{d}$	$k = T^{1/4}$	0.046	0.042	0.046	0.044	0.044	0.046	0.044	0.042	0.043
	$k = T^{1/3}$	0.047	0.042	0.049	0.043	0.444	0.049	0.048	0.040	0.043
	$k = T^{1/2}$	0.052	0.043	0.049	0.044	0.049	0.050	0.049	0.045	0.045

**Table 3.** Estimation of  $\lambda = (d, \phi_1)'$ .

DGP: $ARFIMA(1, d, 0)$ , $\phi_1 = 0.6$ , $k = T^{1/4}$									
$\mu_0$ <b>known</b> ( $\mu_0 = 0$ )									
$d_0$	-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$									
bias $\hat{d}$	-0.030	-0.018	-0.044	0.025	-0.046	-0.053	-0.053	-0.034	-0.055
SRMSE $\hat{d}$	0.212	0.225	0.240	0.230	0.253	0.234	0.237	0.245	0.255
bias $\hat{\phi}_1$	-0.013	-0.023	-0.062	0.000	0.006	0.007	0.011	0.021	0.008
SRMSE $\hat{\phi}_1$	0.219	0.249	0.253	0.203	0.203	0.202	0.193	0.215	0.209
$T = 400$									
bias $\hat{d}$	-0.028	0.028	0.012	0.023	0.024	0.053	0.027	0.022	0.054
SRMSE $\hat{d}$	0.150	0.140	0.136	0.158	0.142	0.164	0.150	0.155	0.162
bias $\hat{\phi}_1$	0.001	0.002	-0.010	0.000	0.003	0.029	0.010	0.007	0.018
SRMSE $\hat{\phi}_1$	0.142	0.135	0.135	0.148	0.131	0.145	0.146	0.141	0.143
$\mu_0$ <b>unknown</b> ( $\mu_0 = 0$ ).									
$d_0$	-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$									
bias $\hat{d}$	-0.022	-0.021	0.000	-0.011	-0.045	-0.021	-0.032	-0.023	-0.022
SRMSE $\hat{d}$	0.219	0.221	0.181	0.219	0.221	0.221	0.210	0.233	0.222
bias $\hat{\phi}_1$	-0.024	-0.043	-0.041	-0.025	0.003	0.000	-0.010	0.010	0.006
SRMSE $\hat{\phi}_1$	0.211	0.197	0.182	0.182	0.216	0.190	0.210	0.215	0.227
$T = 400$									
bias $\hat{d}$	-0.004	-0.031	-0.018	-0.024	-0.023	-0.013	-0.015	-0.010	-0.017
SRMSE $\hat{d}$	0.146	0.141	0.131	0.143	0.140	0.143	0.143	0.144	0.147
bias $\hat{\phi}_1$	0.018	0.004	0.013	0.001	0.003	0.026	0.013	-0.016	0.009
SRMSE $\hat{\phi}_1$	0.143	0.136	0.127	0.147	0.132	0.139	0.138	0.140	0.141

**Table 4.** Estimation of  $\lambda = (d, \theta_1)'$ .

DGP: ARFIMA(0, d, 1), $\theta_1 = 0.5$ . $k = T^{1/4}$									
$\mu_0$ known ( $\mu_0 = 0$ )									
$d_0$	-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$									
bias $\hat{d}$	-0.012	-0.016	-0.010	-0.013	-0.018	-0.012	0.0206	0.004	0.025
SRMSE $\hat{d}$	0.148	0.131	0.148	0.147	0.138	0.164	0.244	0.213	0.260
bias $\hat{\theta}_1$	-0.000	0.002	-0.003	-0.003	0.008	-0.003	-0.084	-0.146	-0.339
SRMSE $\hat{\theta}_1$	0.146	0.127	0.138	0.141	0.131	0.141	0.2158	0.255	0.435
$T = 400$									
bias $\hat{d}$	-0.000	-0.007	-0.007	-0.005	-0.0082	-0.012	-0.0037	-0.009	-0.012
SRMSE $\hat{d}$	0.056	0.058	0.057	0.0574	0.056	0.0552	0.0567	0.0591	0.077
bias $\hat{\theta}_1$	0.00	0.005	0.0029	-0.003	0.0045	0.0036	-0.025	-0.04700	-0.1834
SRMSE $\hat{\theta}_1$	0.056	0.0542	0.0590	0.0582	0.0567	0.0541	0.0699	0.0974	0.2453
$\mu_0$ unknown ( $\mu_0 = 0$ )									
$d_0$	-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$									
bias $\hat{d}$	-0.019	-0.021	-0.008	-0.016	-0.021	-0.018	-0.023	-0.022	-0.0126
SRMSE $\hat{d}$	0.135	0.129	0.147	0.145	0.140	0.148	0.139	0.146	0.146
bias $\hat{\theta}_1$	0.002	0.007	-0.005	-0.000	0.012	0.002	0.008	0.007	-0.006
SRMSE $\hat{\theta}_1$	0.129	0.125	0.135	0.136	0.134	0.135	0.122	0.139	0.142
$T = 400$									
bias $\hat{d}$	-0.004	-0.031	-0.018	-0.024	-0.023	-0.013	-0.015	-0.010	-0.017
SRMSE $\hat{d}$	0.146	0.141	0.131	0.143	0.140	0.143	0.143	0.144	0.147
bias $\hat{\theta}_1$	0.018	0.004	0.013	0.001	0.003	0.026	0.013	-0.016	0.009
SRMSE $\hat{\theta}_1$	0.143	0.136	0.127	0.147	0.132	0.139	0.138	0.140	0.141

**Table 5. Estimation of  $\lambda = (d, \phi_1, \phi_2)'$ .**  
DGP: **ARFIMA(2, d, 0)**,  $\phi_1 = \mathbf{0.6}$ ,  $\phi_2 = \mathbf{-0.65}$ ;  $k = T^{1/4}$

	$\mu_0$ known ( $\mu_0 = 0$ )								
$d_0$	-0.7	-0.3	0.4	0.8	1.0	1.4	1.8	2.0	2.4
$T = 100$									
bias $\hat{d}$	-0.003	-0.031	-0.030	-0.021	-0.022	-0.064	-0.113	-0.168	-0.112
SRMSE $\hat{d}$	0.224	0.228	0.230	0.229	0.223	0.225	0.244	0.281	0.282
bias $\hat{\phi}_1$	0.027	0.074	0.052	0.031	0.041	0.058	0.043	0.049	0.051
SRMSE $\hat{\phi}_1$	0.206	0.219	0.232	0.216	0.226	0.230	0.261	0.302	0.306
bias $\hat{\phi}_2$	-0.059	-0.050	-0.083	-0.069	-0.072	-0.067	-0.053	-0.031	-0.038
SRMSE $\hat{\phi}_2$	0.150	0.171	0.174	0.159	0.165	0.165	0.202	0.371	0.382
$T = 400$									
bias ( $\hat{d}$ )	0.004	-0.011	-0.010	-0.006	-0.012	-0.015	-0.032	-0.051	-0.112
SRMSE $\hat{d}$	0.067	0.063	0.061	0.063	0.064	0.064	0.074	0.087	0.149
bias $\hat{\phi}_1$	-0.003	0.003	0.001	0.001	0.003	0.002	-0.005	-0.019	-0.063
SRMSE $\hat{\phi}_1$	0.054	0.054	0.049	0.051	0.053	0.053	0.059	0.067	0.110
bias $\hat{\phi}_2$	0.003	0.003	-0.003	-0.003	-0.003	0.001	0.012	0.024	0.089
SRMSE $\hat{\phi}_2$	0.047	0.049	0.048	0.049	0.047	0.048	0.052	0.061	0.126

**Table 6. SRMSE for various estimation techniques**

$T = 100$ ; DGP: $\Delta^{d_0} y_t = \varepsilon_t$ , $\mu_0$ known							
$d_0$	-0.3	0.2	0.4	0.8	1.0	1.4	1.8
<i>VR</i>	0.211	0.208	0.193	0.172	0.155	0.154	0.160
<i>BER</i>	0.084	0.083	0.089	0.091	0.086	0.092	0.095
<i>TSB</i>	0.189	0.094	0.109	-	-	-	-
<i>GZW</i>	0.082	0.089	0.096	-	-	-	-
<i>SOW</i>	0.089	0.092	0.069	-	-	-	-

## 5. RESIDUAL-BASED STATISTICS FOR DIAGNOSING CHECKING.

Although parametric estimates present in general better properties than semiparametric ones, these good properties rely heavily on the correct specification of the parametric model. For this reason, a formal test to check the adequacy of the proposed model is very often carried out. A common way of testing it is by checking the assumption of white noise residuals (see Milhoj (1981) and Chen and Deo (2000) for other approaches that do not require the computation of residuals from the fitted model). Box and Pierce (1970) introduced a goodness-of-fit procedure that tests for significant residual autocorrelations. The

Box-Pierce (BP) statistic is defined as

$$Q(k) = T \sum_{i=1}^k \hat{\rho}_{e(\hat{\lambda})}^2(i). \quad (17)$$

They proved that, in the context of ARIMA processes, this statistic is asymptotically  $\chi^2$ -distributed with  $k - p - q$  degrees of freedom, for large  $k$ . Ljung and Box (1978) introduced a modification to the  $Q(k)$ -statistic that improves the approximation to the  $\chi_{k-p-q}^2$  distribution. This is defined as

$$\tilde{Q}(k) = T(T+2) \sum_{i=1}^k (T-k)^{-1} \hat{\rho}_{e(\hat{\lambda})}^2(i).$$

Hong (1996) proposed a generalization of the BP test given by

$$H_T = \left( T \sum_{i=1}^{k(T)} h^2(i/k) \hat{\rho}_{e(\hat{\lambda})}^2(i) - C_T(h) \right) / 2D_T(h)^{1/2},$$

where  $h(\cdot)$  is a suitable chosen kernel and  $k(T)$  verifies that  $\lim_{T \rightarrow \infty} k(T) = \infty$  and  $\lim_{T \rightarrow \infty} k(T)/T \rightarrow 0$  (for the definition of  $C_T$  and  $D_T$  see Hong, 1996). When  $h(\cdot)$  is the truncated kernel, i.e.,  $h(z) = 1$  for  $|z| \leq 1$  and 0 for  $|z| > 1$ , it is obtained

$$H_T^* = \left( T \sum_{i=1}^{k(T)} \hat{\rho}_{e(\hat{\lambda})}^2(i) - k(T) \right) / (2k(T))^{1/2}, \quad (18)$$

a generalization of BP's test when  $k(T) \rightarrow \infty$ . Hong (1996) establishes the asymptotic normality of  $H_T$  for AR models when  $k(T) \rightarrow \infty$  and  $k(T)/T \rightarrow 0$ .

The estimation procedure described in Section 3 provides an immediate check on the adequacy of the specified model, since the criterion function defined in (12), evaluated at the estimated values and multiplied by the sample size, coincides with the BP goodness-of-fit statistic in (17). Due to its simplicity, it is worth analyzing the behavior of this statistic. In the context of stationary ARFIMA processes with known mean, the asymptotic distribution of residual autocorrelations have been examined by Li and McLeod (1986). The following theorem is an extension of Li and McLeod's result to the case where the process is allowed to be nonstationary and to have an unknown mean.

**Theorem 6** *Let  $\hat{\rho}_{ke(\hat{\lambda})}$  be the vector containing the autocorrelations up to lag  $k$  of the residuals, defined in (4) or (6), such that:*

$$\hat{\rho}_{ke(\hat{\lambda})} = \left( \hat{\rho}_{e(\hat{\lambda})}(1), \dots, \hat{\rho}_{e(\hat{\lambda})}(k) \right)' \quad (19)$$

*where  $\hat{\rho}_{e(\hat{\lambda})}(i)$  is defined as in (9). Then, for any fixed  $k$ ,  $\sqrt{T} \hat{\rho}_{ke(\hat{\lambda})}$  is asymptotically normally distributed, with zero mean and variance-covariance matrix given by:*

$$\Sigma = I_k - J_k(\lambda_0) (J_k'(\lambda_0) J_k(\lambda_0))^{-1} J_k'(\lambda_0)$$

where  $J_k(\lambda_0)$  is the limit as  $T$  tends to infinity of the Jacobian matrix of  $\hat{\rho}_e(\lambda_0)$ .

Applying standard results, it is easily seen that  $\Sigma$  is approximately idempotent with rank  $k - p - q - 1$ , for  $k$  large enough. Hence, both  $Q(k)$  and  $\tilde{Q}(k)$  are approximately  $\chi^2$ -distributed with  $k - p - q - 1$  degrees of freedom for large  $T$ .

The previous results imply that the minimum value of the criterion function (12),  $V_{ke}(\hat{\lambda})$ , can be used to test the adequacy of the specified model. Under the null hypothesis of correct specification, the above-mentioned value multiplied by the sample size is distributed approximately  $\chi_{k-p-q-1}^2$ , for large  $T$ , where  $k$  is the number of autocorrelations considered in the criterion function. It is also straightforward to compute the Hong's statistic in (18) from this value, just by multiplying it by the sample size, subtracting the number of included autocorrelations and dividing by the squared root of  $2k(T)$ . Although the asymptotic properties of this statistic remain unknown where long memory processes are considered and their derivation goes beyond the scope of this paper, the following section explores by simulation its finite sample behavior. It also discusses the finite sample properties of both  $Q(k)$  and  $\tilde{Q}(k)$  statistics.

## 6. BEHAVIOR OF THE GOODNESS-OF-FIT TESTS IN FINITE SAMPLES

To evaluate the performance in terms of size and power of the goodness-of-fit tests examined in Section 5, the following experiments have been carried out. First, processes of the form  $\Delta^{d_0} y_t = \varepsilon_t$ ,  $\varepsilon_t \sim NID(0, 1)$  were generated for different values of  $d_0$ . This parameter was estimated in accordance with the method presented in Section 3 and the  $Q(k)$  statistic was computed using the corresponding residuals. The value  $k$  was set equal to 3, 4 and 5 for sample sizes  $T = 150, 400$  and  $500$ , respectively (that is,  $k \sim T^{1/4}$ ). Empirical size at the 5% signification level is calculated using the  $\chi_{k-1, 0.95}^2$  value. Since the Ljung-Box statistic improves the approximation to the  $\chi^2$  distribution, it is usually preferred to the Box-Pierce statistic in applications. Therefore, the values of the  $\tilde{Q}(k)$  are also computed in order to compare the behavior of both tests. Table 7 reports the empirical size of both the Box-Pierce (BP) and Ljung-Box (LB) tests. In agreement with the findings of Li and McLeod (1986), the empirical size is close to the nominal size in both tests. Although the approximation to the  $\chi_{k-1}^2$  distribution is slightly better for the LB test, the difference between both tests decreases as  $T$  increases.

With respect to the power of the test, it will obviously depend on how close is the DGP to the null hypothesis. An ARFIMA(1,  $d_0$ , 0) is been chosen as the true DGP with a value of the autoregressive parameter  $\phi_1$  equal to 0.5, for different values of  $d_0$ . An ARFIMA(0,  $d$ , 0) was estimated instead. Table 8 reports the power of both the Box-Pierce (BP) and Ljung-Box (LB) tests at the 5% nominal signification level. Therefore power has been calculated

with the asymptotic critical values and it has not been size-adjusted. Since both tests are in general undersized, better results would have been achieved after size adjustment. It is seen both tests perform quite similarly. For small sample sizes ( $T = 150$ ) the power is similar to that obtained by other methods proposed in the literature (see, for instance, Delgado and Hidalgo, (1999)). It improves considerably when larger samples sizes are considered, providing good results for sample sizes around 400 or 500.

**Table 7.** Empirical Size of BP and LB. ( $S.L. : 5\%$ )

$d_0$	-0.7	-0.3	0	0.4	0.8	1.4
$T = 150$						
<i>BP</i>	3.5%	3.2%	3.8%	4.0%	3.6%	3.8%
<i>LB</i>	4.5%	4.0%	4.5%	4.1%	4.0%	4.5%
$T = 400$						
<i>BP</i>	6.0%	4.2%	4.8%	4.8%	5.8%	4.5%
<i>LB</i>	6.0%	5.8%	3.5%	5.5%	6.0%	5.0%
$T = 500$						
<i>BP</i>	5.5%	5.0%	5.0%	5.0%	4.5%	4.5%
<i>LB</i>	6.0%	5.0%	5.0%	5.5%	5.0%	4.2%

**Table 8.** Power of BP and LB. ( $S.L.: 5\%$ )

DGP: ARFIMA(1, $d$ , 0)						
$d_0$	-0.7	-0.3	0	0.4	0.8	1.4
$T = 150$						
<i>BP</i>	24.8%	24.2%	24.4%	23.0%	23.8%	21.8%
<i>LB</i>	25.3%	24.4%	27.0%	25.2%	25.4%	21.8%
$T = 400$						
<i>BP</i>	66.6%	68.6%	66.9%	64.4%	68.9%	64.5%
<i>LB</i>	67.3%	69.2%	67.9%	65.6%	70.8%	65.3%
$T = 500$						
<i>BP</i>	81.8%	81.2%	79.8%	82.2%	82.4%	82.8%
<i>LB</i>	81.6%	81.6%	80.0%	83.0%	82.2%	83.4%

Tables 9 and 10 report the results of analogous simulations based on Hong's statistic defined in (18). To compute the size, critical values from a  $N(0, 1)$  distribution have been employed. Since in this case  $k$  is allowed to go to infinity, different values of  $k$  have been employed, namely  $k(T) = T^{1/4}$ ,  $T^{1/3}$  and  $T^{1/2}$ . It is seen that the empirical size is lower than the nominal one but the approximation improves for faster  $k(T)$  and larger sample sizes. Size-corrected power is been reported in order to facilitate comparison among different values of  $k$ . Large sample sizes, ( $T = 400, 500$ ), are needed to obtain a reasonable power.

Also, it can be observed that a slower  $k(T)$  provides better power, result which agrees with Hong's (1996) findings.

**Table 9.** Empirical Size of  $H_T^*$ . (S.L.: 5%)

$d_0$	-0.7	-0.3	0	0.4	0.8	1.4
$T = 150$						
$k=T^{1/4}$	1.2%	1.5%	1.4%	1.1%	1.2%	1.2%
$k=T^{1/3}$	2.2%	2.2%	2.3%	1.8%	1.6%	1.5%
$k=T^{1/2}$	2.5%	2.5%	3.4%	2.8%	2.8%	2.6%
$T = 400$						
$k=T^{1/4}$	1.8%	1.9%	1.5%	2.2%	2.4%	2.1%
$k=T^{1/3}$	2.9%	3.2%	2.1%	2.6%	2.9%	1.5%
$k=T^{1/2}$	5.4%	3.9%	3.1%	3.6%	3.9%	3.2%
$T = 500$						
$k=T^{1/4}$	2.7%	2.1%	2.0%	3.1%	2.9%	2.6%
$k=T^{1/3}$	2.4%	2.3%	2.6%	2.5%	2.6%	3.2%
$k=T^{1/2}$	3.9%	4.2%	3.5%	4.0%	4.5%	5.2%

**Table 10.** Power of  $H_T^*$  (size corrected).

DGP: ARFIMA(1, $d_0$ , 0)						
$d_0$	-0.7	-0.3	0	0.4	0.8	1.4
$T = 150$						
$k=T^{1/4}$	22.3%	16.2%	19.8%	21.2%	20.6%	19.3%
$k=T^{1/3}$	20.7%	14.7%	14.2%	14.2%	16.1%	16.1%
$k=T^{1/2}$	12.1%	10.4%	9.2%	10.1%	13.8%	13.5%
$T = 400$						
$k=T^{1/4}$	54.2%	28.7%	65.4%	54.3%	54.9%	56.5%
$k=T^{1/3}$	55.9%	56.4%	65.3%	56.4%	56.5%	59.8%
$k=T^{1/2}$	30.4%	32.6%	39.3%	35.0%	35.5%	30.6%
$T = 500$						
$k=T^{1/4}$	70.9%	82.1%	76.7%	75.0%	74.1%	76.8%
$k=T^{1/3}$	75.9%	77.7%	68.6%	76.6%	77.1%	69.9%
$k=T^{1/2}$	45.2%	52.8%	45.8%	46.1%	43.2%	39.7%

## 7. EMPIRICAL APPLICATIONS.

In order to illustrate the application of the techniques proposed in this paper, two empirical studies have been carried out. Firstly, to provide a further comparison with previous estimation techniques, the empirical series analyzed by Beran (1995), Velasco and Robinson

(2000) and Robinson (1994b) have been considered, namely the chemical process temperature readings (Series C) and the chemical process concentration readings (Series A) from Box and Jenkins (1976). Secondly, the orders of fractional integration of the *GDP* per capita series of several countries have been estimated.

With respect to the first application, our conclusions are in fair agreement with those of the above-mentioned studies. For series C, Box and Jenkins (BJ), fitted an ARIMA(1,1,0) model, with an estimated value of the AR parameter  $\phi_1 = 0.8$ . The corresponding estimates for an ARFIMA(1,  $d$ , 0) process according with the GMD method are:  $\hat{d} = 1.005$  with a 95% confidence interval (C.I.) of [0.7497, 1.2617] and  $\hat{\phi}_1$  is 0.798 with a 95% C.I. of [0.563, 0.972], in close agreement with BJ's conclusions. Nevertheless, the recognition of the uncertainty on  $d$ , increases substantially the standard deviation of the AR parameter. Similar conclusions are reached both in Beran (1995) and Velasco and Robinson (2000).

For series A, BJ fitted an ARIMA(0,1,1), which yields a significant value for the MA parameter equal to -0.7. This large negative value suggests that the model could be overdifferenced. If an ARFIMA(0,  $d$ , 1) is fitted instead, the GMD estimates are 0.43 and -0.038 with 95% C.I.s of [0.241,0.612] and [-0.27,0.196] for  $\hat{d}$  and  $\hat{\theta}_1$ , respectively. Therefore,  $d = 1$  is not included in the C.I., which reinforces the believe that the series in the BJ model is overdifferenced. Since the MA parameter is not significant in this second case, we have also fitted an ARFIMA(0,  $d$ , 0) to the data. In this case, the estimated value of  $d$  drops to 0.401 with a 95% C.I. of [0.292,0.51]. Also in this case, our results follow closely the ones obtained in Velasco and Robinson (2000) and Beran (1995).

In the second application, some of the series of Maddison's (1995) data set have been analyzed. This data set contains the annual *GDP* per capita series for OECD countries during the period 1870-1994 (125 observations). This data set has also been analyzed by Michelaci and Zaffaroni (2000) and Dolado et al. (2002) among others. In particular, three countries have been considered, namely Canada, Japan and Germany. Since these series are clearly trended, the estimation has been carried out as if the mean was different from zero and unknown. They have been estimated according to the following parametric methods: the GMD method presented in this paper, the Sowell's exact ML procedure (with the series in first differences when they are nonstationary), the NLS method by Beran (1995) and the Velasco and Robinson (2000) Whittle estimator (with Zhurbenko taper of order 2). Also, several ARFIMA( $p, d, q$ ) processes have estimated for values of  $p, q$  in the range 0 to 2, and the results reported correspond to the preferred model according to the AIC lag-length criterion. Table 11 reports the value of the estimates and their standard errors (in brackets). As it can be seen from this table, all methods deliver very similar estimates of the parameters. With the exception of Japan, it seems to be the case that Canada and Germany are clear cases where the *GDP* per capita series are fractionally integrated, with

a value of  $d$  around 0.5 for the former and in the range  $d \in (0.5, 1)$  for the latter. It is also noticeable that there is a negative relation between the values of  $d$  and the values of the remaining parameters across the different methods. Standard deviations are also similar across methods with the exception of the Velasco and Robinson technique which shows slightly higher values due to the tapering employed.

**Table 11.** Estimation results from various estimation techniques

	Canada	Japan	Germany
Model (AIC)	ARFIMA(1, $d$ , 0)	ARFIMA(0, $d$ , 0)	ARFIMA(0, $d$ , 1)
GMD	$\hat{d} = 0.40; \hat{\phi}_1 = 0.78$ (0.217; 0.174)	$\hat{d} = 1.083$ (0.006)	$\hat{d} = 0.87; \hat{\theta}_1 = 0.49$ (0.195; 0.218)
SOW	$\hat{d} = 0.48; \hat{\phi}_1 = 0.72$ (0.239; 0.213)	$\hat{d} = 1.064$ (0.006)	$\hat{d} = 0.80; \hat{\theta}_1 = 0.51$ (0.201; 0.222)
BER	$\hat{d} = 0.50; \hat{\phi}_1 = 0.67$ (0.242; 0.230)	$\hat{d} = 1.067$ (0.006)	$\hat{d} = 0.78; \hat{\theta}_1 = 0.55$ (0.214; 0.229)
V-R	$\hat{d} = 0.41; \hat{\phi}_1 = 0.64$ (0.345; 0.340)	1.030 (0.101)	$\hat{d} = 0.62; \hat{\theta}_1 = 0.62$ (0.330; 0.341)

## 8. CONCLUSIONS

In this paper we have proposed a new method for estimating the parameters of an ARFIMA( $p, d_0, q$ ) process with  $d_0 > -0.75$ . It covers a very wide range of values of  $d_0$ , providing therefore a unified framework for the construction of confidence intervals and tests for the memory parameter. The proposed estimator belongs to the MD class and it is based on the minimization of the residuals obtained after filtering a process through ARFIMA parameters. Its asymptotic properties as well as its finite sample performance are discussed and it is shown that it is  $\sqrt{T}$ -consistent, asymptotically normally distributed, and efficient without making strong assumptions on the distribution of the process under study. Monte Carlo experiments show that it is also well-behaved in finite samples and that it compares well to other existing estimators in the literature. Another interesting feature of the estimator is that the criterion function evaluated in the estimate coincides with the Box-Pierce (1970) goodness-of-fit statistic, providing therefore, and immediate tool to evaluate the adequacy of the model specification. The asymptotic properties of this statistic, as well as the ones of the Ljung-Box (1978) statistic, are discussed and some simulations are provided in order to evaluate their accuracy in finite samples.

Finally, another nice attribute of the proposed estimator is its flexibility to be extended to more general settings. For instance, the estimator can be easily robustified against conditional heterocedasticity, simply by considering the sample autocorrelations of the standardized residuals and, with this modification, the resulting statistic remains asymptotically equivalent to MLE. Also, following the lines of Wright (1999), it can be adapted to deal with the fractionally integrated stochastic volatility model. Further research should also be

undertaken to extend the previous framework to the multivariate case.

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## APPENDIX 1

### Proof of Theorem 1.

The residuals defined in (6),  $e_t(\lambda)$ , evaluated at  $\lambda = \lambda_0$  can be written as

$$e_t(\lambda_0) = \sum_{j=0}^{t-m_0-1} \alpha_j(\lambda_0^*) (\Delta^{m_0} y_t - \mu_0) + (\overline{\Delta^{m_0} y} - \mu_0) \sum_{j=0}^{t-m_0-1} \alpha_j(\lambda_0^*), \quad t = m_0 + 1, \dots, T.$$

This allows us to write

$$e_t(\lambda_0) = \varepsilon_t + \eta_t,$$

where  $\eta_t$  is given by

$$\eta_t = - \sum_{i=t-m_0}^{\infty} \alpha_i(\lambda_0^*) (\Delta^{m_0} y_{t-i} - \mu_0) + (\overline{\Delta^{m_0} y} - \mu_0) \sum_{j=0}^{t-m_0-1} \alpha_j(\lambda_0^*),$$

and  $\overline{\Delta^{m_0} y}$  denotes the sample mean of  $\Delta^{m_0} y_t$ . To simplify notation, assume without loss of generality that  $m_0 = 0$  (when this is not the case, define the process  $y_t^* = \Delta^{m_0} y_t$  and the following arguments will remain valid just by substituting  $y_t$  by  $y_t^*$ ). Then,  $d_0 = \varphi_0$ .

Let  $\sqrt{T}\Upsilon_T$  be the vector that contains the differences in expression (11), that is

$$\sqrt{T}\Upsilon_T = \sqrt{T} \left( \hat{\rho}_{ke(\lambda_0)} - \hat{\rho}_{k\varepsilon} \right).$$

Consider the  $i$ -th element of the vector  $\Upsilon_T$ . The sample variance of  $e_t(\lambda_0)$  converges to the innovation variance as long as  $d_0 > -1$  (see Odaki, 1993). The difference between the  $i$ -th sample autocovariance of the residuals and of the  $i$ -th autocovariance of the innovations, scaled by  $\sqrt{T}$ , is given by

$$T^{1/2} \left( \frac{\sum_{t=1}^{T-i} e_t(\lambda_0) e_{t+i}(\lambda_0) - \sum_{t=1}^{T-i} \varepsilon_t \varepsilon_{t+i}}{T} \right) = \quad (20)$$

$$= T^{-1/2} \left( \sum_{t=1}^{T-i} \varepsilon_{t+i} \eta_t + \sum_{t=1}^{T-i} \varepsilon_t \eta_{t+i} + \sum_{t=1}^{T-i} \eta_t \eta_{t+i} \right). \quad (21)$$

Let us first consider the case where  $\mu_0$  is known and equal to zero. In this case  $\eta_t$  collapses to  $\eta_t = - \sum_{i=t}^{\infty} \alpha_i(\lambda_0^*) y_{t-i}$ . By repeated substitution, Odaki (1993) shows that this expression can be rewritten as

$$\eta_t = \sum_{j=0}^{\infty} \psi_{j,t-1} \varepsilon_{-j}, \quad (22)$$

(for the precise form of the sequence of coefficients  $\{\psi_{j,t-1}\}_{j=0}^{\infty}$ , see Odaki, 1993, p. 704). Notice that only terms of  $\varepsilon_t$  with  $t \leq 0$  enter the above definition. From the former article it is also known that the orders of magnitude of the sum of squares of these coefficients are

$$\left( \sum_{j=0}^{\infty} \psi_{j,t-1}^2 \right) = \begin{cases} O(t^{-1}) & \text{if } \varphi_0 \in (-0.5, 0.5), \\ O((\log t)t^{-1}) & \text{if } \varphi_0 = -0.5, \\ O(t^{-2(1+\varphi_0)}) & \text{if } \varphi_0 < -0.5. \end{cases} \quad (23)$$

To check that expression (21) converges to zero, consider

$$T^{-1/2} \sum_{t=1}^{T-i} E(\varepsilon_{t+i}\eta_t) = T^{-1/2} \sum_{t=1}^{T-i} E(\varepsilon_{t+i}) E(\eta_t) = 0, \quad (24)$$

by independence of the processes  $\eta_t$  and  $\{\varepsilon_t\}_{t=1}^{\infty}$  (see (22)). Taking into account (24) and (23)<sup>3</sup>,

$$T^{-1} E \left( \sum_{t=1}^{T-i} \varepsilon_{t+i}^2 \eta_t^2 \right) = \sigma^4 T^{-1} \sum_{t=1}^{T-i} \left( \sum_{j=0}^{\infty} \psi_{j,t-1}^2 \right) \rightarrow 0. \quad (25)$$

A similar argument holds for the second term in (21), since in this case  $\eta_{t+i} = \sum_{j=0}^{\infty} \psi_{j,t+i-1} \varepsilon_{-j}$  and again no contemporaneous terms of  $\varepsilon$  are found in the product  $\varepsilon_t \eta_{t+i}$ . Hence, all cross products vanish. With respect to the third term in (21),

$$\begin{aligned} T^{-1/2} \sum_{t=1}^{T-i} E |\eta_t \eta_{t+i}| &\leq T^{-1/2} \sum_{t=1}^{T-i} E(\eta_t^2) \\ &= T^{-1/2} \sigma^2 \sum_{t=1}^{T-i} \left( \sum_{j=0}^{\infty} \psi_{j,t-1}^2 \right). \end{aligned} \quad (26)$$

The orders of magnitude in (23) imply that

$$\sum_{t=1}^{T-i} \left( \sum_{j=0}^{\infty} \psi_{j,t-1}^2 \right) = \begin{cases} O(\log T) & \text{for } \varphi_0 \in (-0.5, 0.5) \\ O(\log^2 T) & \text{for } \varphi_0 = -0.5 \\ O(T^{-2(1+\varphi_0)+1}) & \text{for } \varphi_0 < -0.5. \end{cases} \quad (27)$$

Therefore, for  $-0.75 < \varphi_0 < 0.5$ , expression (26) tends to zero. Note that if we consider the covariances (without being multiplied by  $\sqrt{T}$ ), expression (26) still tends to zero even in the interval  $-1 < \varphi_0 < -0.75$ . To check that the variance also converges to zero, notice that

$$T^{-1} E \left( \sum_{t=1}^{T-i} \eta_t \eta_{t+i} \right)^2 = T^{-1} \sum_{r=1}^{T-i} \sum_{s=1}^{T-i} E(\eta_r \eta_{r+i} \eta_s \eta_{s+i}), \quad (28)$$

and that

$$E(\eta_r \eta_{r+i} \eta_s \eta_{s+i}) \leq E(\varepsilon_t^4) \sum_{j=0}^{\infty} |\psi_{j,r-1} \psi_{j,r+i-1} \psi_{j,s-1} \psi_{j,s+i-1}| +$$

---

<sup>3</sup>Since  $\sum_{j=0}^{\infty} E(\psi_{j,t-1}^2 \varepsilon_{-j}^2) < \infty$ , it is possible to interchange the expectation and the summation in order to obtain (25). See Rao, (1973), p.111.

$$\begin{aligned}
& +\sigma^4 \sum_{j=0}^{\infty} |\psi_{j,r-1}\psi_{j,s-1}| \sum_{j=0}^{\infty} |\psi_{j,r+i-1}\psi_{j,s+i-1}| \\
& +\sigma^4 \sum_{j=0}^{\infty} |\psi_{j,r-1}\psi_{j,s+i-1}| \sum_{j=0}^{\infty} |\psi_{j,r+i-1}\psi_{j,s-1}| \\
& +\sigma^4 \sum_{j=0}^{\infty} |\psi_{j,r-1}\psi_{j,r+i-1}| \sum_{j=0}^{\infty} |\psi_{j,s-1}\psi_{j,s+i-1}|.
\end{aligned}$$

It is easy to check that the coefficients  $\{\psi_{j,t}\}$  are strictly smaller than one in absolute value and that are strictly decreasing in both subindexes  $(j, t)$ . Applying the last two results and Cauchy's inequality, we obtain:

$$\begin{aligned}
E(\eta_r \eta_{r+i} \eta_s \eta_{s+i}) & \leq E(\varepsilon_t^4) \left\{ \left( \sum_{j=0}^{\infty} |\psi_{j,r-1}^2 \psi_{j,r+i-1}^2| \sum_{j=0}^{\infty} |\psi_{j,s-1}^2 \psi_{j,s+i-1}^2| \right)^{1/2} \right. \\
& \quad \left. + 2 \left( \sum_{j=0}^{\infty} |\psi_{j,r-1} \psi_{j,s-1}| \right)^2 + \sum_{j=0}^{\infty} |\psi_{j,r-1}^2| \sum_{j=0}^{\infty} |\psi_{j,s-1}^2| \right\} \\
& \leq 4E(\varepsilon_t^4) \sum_{j=0}^{\infty} |\psi_{j,r-1}^2| \sum_{j=0}^{\infty} |\psi_{j,s-1}^2|.
\end{aligned}$$

Therefore

$$T^{-1} \sum_{r=1}^{T-i} \sum_{s=1}^{T-i} E(\eta_r \eta_{r+i} \eta_s \eta_{s+i}) \leq 4T^{-1} E(\varepsilon_t^4) \sum_{r=1}^{T-i} \sum_{s=1}^{T-i} \left( \sum_{j=0}^{\infty} \psi_{j,r-1}^2 \right) \left( \sum_{j=0}^{\infty} \psi_{j,s-1}^2 \right), \quad (29)$$

and again taking into account (27), this expression tends to zero for  $\varphi_0 > -0.75$ .

Since the  $i$ -th element of  $\sqrt{T}\Upsilon_T$  tends to zero for all  $i = 1, \dots, k$ , then  $\sqrt{T}\Upsilon_T \xrightarrow{p} 0$ , implying the desired result.

The case where  $\mu_0$  is unknown can be proved similarly using standard arguments since  $\overline{\Delta}^{m_0 y}$  is a consistent estimator of  $\mu_0$  (see for instance, Robinson 1994b).■

## Proof of Theorem 2

Let us first consider the convergence of  $\sqrt{T}A_{k(T)}\hat{\rho}_{k(T)e(\lambda_0)} \xrightarrow{w} Z$ , where  $Z \sim N_h(0, \cdot)$ . Recall that  $\hat{\rho}_{k(T)e(\lambda_0)} = \hat{\sigma}^{-1}\hat{\gamma}_{k(T)e(\lambda_0)}$ , where  $\hat{\gamma}_{k(T)e(\lambda_0)}$  is the autocovariance function, and  $\hat{\sigma}$  is a consistent estimator of the variance of  $\varepsilon_t$ . We would first show that  $\sqrt{T}A_{k(T)}\hat{\gamma}_{k(T)e(\lambda_0)} \xrightarrow{w} Z^* = \sigma Z$ , from where the corresponding result for  $\hat{\rho}_{k(T)e(\lambda_0)}$  follows trivially.

Following Proposition 6.3.9 in Brockwell and Davis (1993), in order to prove the former statement, it is needed to show that: *i*) for finite  $r$ ,  $\sqrt{T}A_r\hat{\gamma}_{re(\lambda_0)} \xrightarrow{w} Z_r^*$ , where  $Z_r^* \sim N_h(0, \sigma^2 A_r A_r')$ , *ii*)  $Z_r^* \xrightarrow{w} Z^*$ , where  $Z^* \sim N_h(0, \sigma^2 \cdot)$  and  $\sigma^2 = \lim_{r \rightarrow \infty} A_r A_r'$ , and *iii*)

$$\lim_{r \rightarrow \infty} \limsup_{T \rightarrow \infty} P \left( \left| \sqrt{T}A_{k(T)}\hat{\gamma}_{k(T)e(\lambda_0)} - \sqrt{T}A_r\hat{\gamma}_{re(\lambda_0)} \right| > \epsilon \right) = 0 \quad (30)$$

for all  $\epsilon > 0$ . The proof of i) follows from the same arguments as Theorem 1, since for finite  $r$ , the vector  $\hat{\gamma}_{r, e(\lambda_0)}$  is jointly normal. As regards ii), it is a standard result that if  $Z_r^* \sim N_h(0, \sigma^2 A_r A_r')$  and  $\lim_{r \rightarrow \infty} A_r A_r' = \Sigma$ , then  $Z_r^* \xrightarrow{w} Z^*$  (see for instance Brockwell and Davis, 1993, p.225). To prove iii), notice that expression (30) can be rewritten as

$$\lim_{r \rightarrow \infty} \limsup_{T \rightarrow \infty} P \left( \left| \sqrt{T} \sum_{j=r+1}^{k(T)} a_{ij} \hat{\gamma}_{e(\lambda_0)}(j) \right| > \epsilon \right) = 0, \quad \text{for } i = 1, \dots, h. \quad (31)$$

By Chebyshev's inequality, the probability in (31) is bounded by

$$\begin{aligned} & \epsilon^{-2} TE \left( \left( \sum_{j=r+1}^{k(T)} a_{ij} \hat{\gamma}_{e(\lambda_0)}(j) \right)^2 \right) \\ &= \epsilon^{-2} \left( \sum_{j=r+1}^{k(T)} a_{ij}^2 TE \left( \hat{\gamma}_{e(\lambda_0)}(j) \right)^2 + 2 \sum_{s=r+1}^{k(T)} \sum_{l>s}^{k(T)} a_{il} a_{is} TE \left( \hat{\gamma}_{e(\lambda_0)}(s) \hat{\gamma}_{e(\lambda_0)}(l) \right) \right) \end{aligned}$$

It is easy to check that  $TE \left( \hat{\gamma}_{e(\lambda_0)}(j) \right)^2 = \frac{(T-j)}{T} \sigma^4$  and that  $TE \left( \hat{\gamma}_{e(\lambda_0)}(s) \hat{\gamma}_{e(\lambda_0)}(l) \right) = 0$ . Then,

$$\lim_{r \rightarrow \infty} \lim_{T \rightarrow \infty} \epsilon^{-2} \left( \sum_{j=r+1}^{k(T)} a_{ij}^2 TE \left( \hat{\gamma}_{e(\lambda_0)}(j) \right)^2 + 2 \sum_{s=r+1}^{k(T)} \sum_{l>s}^{k(T)} a_{il} a_{is} TE \left( \hat{\gamma}_{e(\lambda_0)}(s) \hat{\gamma}_{e(\lambda_0)}(l) \right) \right) \quad (32)$$

$$= \lim_{r \rightarrow \infty} \lim_{T \rightarrow \infty} \epsilon^{-2} \sum_{j=r+1}^{k(T)} a_{ij}^2 \frac{(T-j)}{T} = \lim_{r \rightarrow \infty} \epsilon^{-2} \sum_{j=r+1}^{\infty} a_{ij}^2 = 0 \quad (33)$$

This last equality follows from noticing that the sum  $\sum_{j=1}^{\infty} a_{ij}^2$ ,  $i = 1, \dots, h$ , is bounded since it corresponds to the diagonal elements of the matrix  $\Sigma$ , which is bounded. Expression (33) implies that (30) tends to zero. Therefore,  $\sqrt{T} A_{k(T)} \hat{\gamma}_{k(T)e(\lambda_0)} \xrightarrow{w} \sigma Z$ . Now, we check that  $\left\| \sqrt{T} \left( \hat{A}_{k(T)} - A_{k(T)} \right) \hat{\gamma}_{k(T)e(\lambda_0)} \right\| = o_p(1)$ .

$$\begin{aligned} \sqrt{T} \left\| \left( \hat{A}_{k(T)} - A_{k(T)} \right) \hat{\gamma}_{k(T)e(\lambda_0)} \right\| &\leq \sqrt{T} \left\| \left( \hat{A}_{k(T)} - A_{k(T)} \right) \right\| \left\| \hat{\gamma}_{k(T)e(\lambda_0)} \right\| \\ &= \left( \sum_{i=1}^p \sum_{j=1}^k (a_{ij} - \hat{a}_{ij})^2 \right)^{1/2} \left( T \sum_{j=1}^k \left( \hat{\gamma}_{je(\lambda_0)} \right)^2 \right)^{1/2} \quad (34) \end{aligned}$$

By applying Jensen's inequality and noticing that  $k/T \rightarrow 0$ , it is obtained that  $(T/k)^{1/2} E \left( \sum_{j=1}^k \left( \hat{\gamma}_{je(\lambda_0)} \right)^2 \right)^{1/2} = O_p(1)$ , and that  $(T/k) \text{var} \left( \sum_{j=1}^k \left( \hat{\gamma}_{je(\lambda_0)} \right)^2 \right)^{1/2} \rightarrow 0$ . Then, the second term in (34) is  $O_p(k^{1/2})$ . The first term is  $o_p(k^{-1/2})$ , since  $\left\| \hat{A}_{k(T)} - A_{k(T)} \right\| = o_p(k^{-1/2})$  by hypothesis. This implies that  $\sqrt{T} \hat{A}_{k(T)} \hat{\gamma}_{k(T)e(\lambda_0)} \xrightarrow{w} \sigma Z$ . Given the consistency of  $\hat{\sigma}$ , it follows that  $\sqrt{T} \hat{A}_{k(T)} \hat{\rho}_{k(T)e(\lambda_0)} \xrightarrow{w} Z$  as desired. ■

In order to prove the consistency of the estimator stated in Theorem 3, we consider separately the cases where the inferior limit of the parameter space of  $d$ ,  $\nabla_1$ , is such that  $\nabla_1 \leq d_0 - 1/2$  from those where  $\nabla_1 > d_0 - 1/2$ . The reason for making this distinction is the non-uniform behavior of  $FI(d_0)$  processes which, in turn, determines the properties of the criterion function: whenever  $d_0$  is  $\geq 1/2$ , the process is not stationary and correlations do not exist. When defining the residuals, the process  $y_t$  will be filtered by  $\Delta^d$ , and the resulting process will be a  $FI(d_0 - d)$  process. If  $\nabla_1 > d_0 - 1/2$ , then  $d_0 - d < 1/2$  always, and then the autocorrelations of the above mentioned residuals will exist. A similar problem and solution appeared in Robinson (1995) and Velasco and Robinson (2000). A two-step proof similar to that proposed in both articles is also followed in this paper. Before stating the proof, we need the following definition.

**Definition 1** Denote  $\lambda^{(1)} = d$  and  $\Lambda_1 = \{d : \nabla_1 \leq d \leq \nabla_2\} \times \Lambda^{(-1)}$  if  $\nabla_1 > d_0 - 1/2$  or otherwise  $\Lambda_1 = \{d : d_0 - 1/2 + \eta \leq d \leq \nabla_2\} \times \Lambda^{(-1)}$  for some  $\eta \in (0, 1/2)$ , if  $\nabla_1 \leq d_0 - 1/2$ , where  $\Lambda^{(-1)}$  is the parameter space of the remaining ARMA parameters.

The following auxiliary result is needed.

**Lemma 1** Let  $V_{ke}(\lambda)$  be the criterion function in (12), where  $e_t(\lambda)$  is defined as in (4) or (6), according to the case where the DGP has a known or unknown mean respectively. Let  $\tilde{\lambda} \in \Lambda_1$  and define  $V_{k\varepsilon}(\tilde{\lambda}) = \sum_{i=1}^k \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2$ , where  $\rho_{\varepsilon(\tilde{\lambda})}(\cdot)$  are the (population) autocorrelations associated to the non truncated residuals  $\varepsilon(\tilde{\lambda})$ , and  $k$  is a fixed number. Then:

1.  $V_{ke}(\lambda)$  is continuous in  $\lambda$ ,  $V_{ke}(\tilde{\lambda})$  converges in probability to  $V_{k\varepsilon}(\tilde{\lambda})$  and the convergence is uniform.
2.  $V_{k\varepsilon}(\tilde{\lambda})$  is a continuous function and has a unique minimum at  $\lambda_0$ , such that  $V_{k\varepsilon}(\lambda_0) = 0$ .

**Proof of Lemma 1.**

1. The continuity of  $V_{ke}(\lambda)$  is trivial, since it is a continuous composition of continuous functions. The asymptotic negligibility of the truncation is proved in Proposition 7.3.3. in Brockwell and Davis (1993). Since the sample correlations associated to stationary processes are consistent (Hosking, 1996), it follows that  $V_{ke}(\tilde{\lambda}) \xrightarrow{p} V_{k\varepsilon}(\tilde{\lambda})$ . The uniform convergence follows from the pointwise convergence and an equicontinuity argument using the compactness of  $\Lambda_1$  and the differentiability of  $\rho(\cdot)$  with respect to  $\lambda$  (cf. Davidson, (1994), p. 340, and Velasco and Robinson, 2000).

2. It is straightforward to check that it has a unique minimum at  $\lambda_0$ , since  $\varepsilon_t(\lambda)|_{\lambda=\lambda_0} = \varepsilon_t$ , which is an *i.i.d.* process and therefore all its correlations are zero, (which implies that  $V_{k\varepsilon}(\lambda_0) = 0$ ), but presents non-null autocorrelations for any other value of  $\tilde{\lambda} \neq \lambda_0$  (and therefore  $V_{k\varepsilon}(\tilde{\lambda}) > 0$ ). The continuity of  $V_{k\varepsilon}(\tilde{\lambda})$  follows from the assumptions above (see Amemiya, Theorem 4.1.1) ■

### Proof of Theorem 3

We proceed with the two-step proof of consistency proposed in Robinson (1995) and Velasco and Robinson (2000).

*First step.* Define  $\hat{\lambda}_1 = \arg \min_{\lambda \in \Lambda_1} V_{k\varepsilon}(\lambda)$ . It follows from standard results that if we can write  $V_{k\varepsilon}(\lambda) - V_{k\varepsilon}(\lambda_0) = S(\lambda) - U(\lambda)$ , where  $S(\lambda)$  is nonstochastic and constant over  $t$  such that for all  $\epsilon > 0$  there exists  $\delta > 0$  such that  $\inf_{\|\lambda - \lambda_0\| \geq \epsilon} S(\lambda) \geq \delta$ , and  $\sup_{\lambda \in \Lambda_1} |U(\lambda)| \xrightarrow{p} 0$ , then  $\hat{\lambda}_1 \xrightarrow{p} \lambda_0$ . Hence, let us denote  $S(\lambda) = V_{k\varepsilon}(\lambda)$  and, since  $V_{k\varepsilon}(\lambda)$  is continuous and has a unique minimum at  $\lambda = \lambda_0$ , the condition on  $S(\lambda)$  holds;  $U(\lambda)$  is given by  $U(\lambda) = V_{k\varepsilon}(\lambda) - V_{k\varepsilon}(\lambda) + V_{k\varepsilon}(\lambda_0)$ . Notice that:

$$\sup_{\lambda \in \Lambda_1} |U(\lambda)| \leq \sup_{\lambda \in \Lambda_1} |V_{k\varepsilon}(\lambda) - V_{k\varepsilon}(\lambda)| + |V_{k\varepsilon}(\lambda_0)|,$$

and both terms in the right hand side tend to zero, the first due to uniform convergence and the second due to pointwise convergence (Lemma 1).

*Second step.* Now consider the case where  $d_0 \geq 1/2 + \nabla_1$  and define  $\Lambda_2 = \{d : \nabla_1 \leq d < d_0 - 1/2 + \eta\} \times \Lambda^{(-1)}$ ,  $\hat{\lambda}_k = \arg \min_{\lambda \in \Lambda} V_{k\varepsilon}(\lambda)$ . Now we prove that the probability of the limit of  $\hat{\lambda}_k$  is different from the limit of  $\hat{\lambda}_1$ , converges to zero. For this, it is needed to show that, for any  $\delta > 0$ ,  $P\left(\left\|\hat{\lambda}_k - \hat{\lambda}_1\right\| \geq \delta\right) \rightarrow 0$ . Notice that,

$$P\left(\left\|\hat{\lambda}_k - \hat{\lambda}_1\right\| \geq \delta\right) \leq P\left(\inf_{\lambda \in \Lambda_2} V_{k\varepsilon}(\lambda) \leq \min_{\lambda \in \Lambda_1} V_{k\varepsilon}(\lambda)\right) \quad (35)$$

$$= P\left(\inf_{\lambda \in \Lambda_2} V_{k\varepsilon}(\lambda) \leq V_{k\varepsilon}(\hat{\lambda}_1)\right) \quad (36)$$

$$= P\left(\inf_{\lambda \in \Lambda_2} V_{k\varepsilon}(\lambda) - V_{k\varepsilon}(\hat{\lambda}_1) \leq 0\right) \quad (37)$$

Since  $\hat{\lambda}_1$  is consistent (step 1) and uniform convergence holds,  $P\left(V_{k\varepsilon}(\hat{\lambda}_1) > 0\right) \rightarrow 0$ . On the other hand,  $P(V_{k\varepsilon}(\lambda) \leq 0) \rightarrow 0$  for all  $\lambda \in \bar{\Lambda}_2$ , where  $\bar{\Lambda}_2$  is the closure of  $\Lambda_2$ . To check that, notice that the function  $V_{k\varepsilon}(\lambda)$ ,  $\lambda \in \bar{\Lambda}_2$ , contains the squared estimated correlations of a  $FI(d_0 - \lambda^{(1)})$  process, where  $\lambda^{(1)} \in [\nabla_1, d_0 - 1/2 + \eta]$ , therefore it is always a non-negative quantity. Whenever  $\lambda^{(1)} \in (d_0 - 1/2, d_0 - 1/2 + \eta]$ , then  $1/2 - \eta < d_0 - \lambda^{(1)} < 1/2$ , and the corresponding filtered process is long-memory stationary. Thus, the squared sample autocorrelations converge to the squared population autocorrelations which, clearly, are

bounded away from zero. If  $\lambda^{(1)} \in [\nabla_1, d_0 - 1/2]$ , then  $d_0 - \lambda^{(1)} \geq 1/2$ , so that  $V_{ke}(\lambda)$  contains the sample autocorrelations of a non-stationary process. Since  $V_{ke}(\lambda)$  should contain, at least, the first autocorrelation, it is clear that  $V_{ke}(\lambda) = \sum_{i=1}^k \hat{\rho}_{e(\lambda)}^2(i) \geq \hat{\rho}_{e(\lambda)}^2(1) \xrightarrow{P} 1$  (see Sowell, (1990), Theorem 3). This implies that  $P(V_{ke}(\lambda) \leq 0) \rightarrow 0$  for all  $\lambda \in \bar{\Lambda}_2$  and. The continuity of  $V_{ke}(\lambda)$  implies that infimum is contained in  $\bar{\Lambda}_2$ , which in turn implies that  $P(\inf_{\lambda \in \Lambda_2} V_{ke}(\lambda) \leq 0) \rightarrow 0$ . The latter result, together with  $P(V_{ke}(\hat{\lambda}_1) > 0) \rightarrow 0$ , involves that the probability in (35) tends to zero. ■

#### Proof of Theorem 4

The mean value theorem applied to the first-order condition gives

$$0 = \frac{\partial V_{ke}(\lambda_0)}{\partial \lambda} + \frac{\partial^2 V_{ke}(\bar{\lambda})}{\partial \lambda \partial \lambda'} (\hat{\lambda}_k - \lambda_0),$$

where  $\bar{\lambda}$  is a mean value on the line joining  $\hat{\lambda}_k$  and  $\lambda_0$ . Multiplying through by  $\sqrt{T}$  and solving for  $\sqrt{T}(\hat{\lambda}_k - \lambda_0)$  yields

$$\sqrt{T}(\hat{\lambda}_k - \lambda_0) = - \left( \frac{\partial \hat{\rho}'_{ke}(\bar{\lambda})}{\partial \lambda} \frac{\partial \hat{\rho}_{ke}(\bar{\lambda})}{\partial \lambda'} \right)^{-1} \frac{\partial \hat{\rho}'_{ke}(\lambda_0)}{\partial \lambda} \sqrt{T} \hat{\rho}_{ke}(\lambda_0). \quad (38)$$

Given that  $\hat{\lambda}_k$  is consistent, so is  $\bar{\lambda}$ . To compute the limit of the first product in the right hand side of (38) define  $\hat{J}_k(\lambda_0) = \partial \hat{\rho}_{ke}(\bar{\lambda}) / \partial \lambda' \big|_{\lambda=\lambda_0}$  and notice that  $e_t(\lambda)$  can be written as

$$e_t(\lambda) = \Delta^{d-d_0} \Phi(L)^{-1} \Theta(L) \Phi_0(L) \Theta_0(L)^{-1} \varepsilon_t, \quad (39)$$

or as,

$$e_t(\lambda) = \Theta^{-1}(L) \Phi(L) \Theta_0(L) \Phi_0^{-1}(L) \Delta^{d-d_0} \varepsilon_t + \Theta^{-1}(L) \Phi(L) \left( \Delta^{d-m_0} \mu - \Delta^\varphi \left( \overline{\Delta^{d-\varphi} y} \right) \right), \quad (40)$$

according to the cases where  $\mu$  is known or unknown respectively. Let us analyze the former case.

Consider first the  $(i, 1)$  element of  $\hat{J}_k(\lambda_0)$ . Noticing that  $\partial \Delta^{d-d_0} / \partial d \big|_{d=d_0} = \log(1-L)$  and using the similar arguments as in proof of Theorem 1, it is obtained that,

$$\partial \hat{\rho}_e(i) / \partial d \big|_{\lambda=\lambda_0} = \frac{1}{T} \frac{\sum_{t=i+1}^T \log(1-L) \varepsilon_t \varepsilon_{t-i}}{\sigma^2} + o_p(1), \text{ for } i = 1, \dots, k, \quad (41)$$

(see Tanaka (1999) for details). Now, since  $\log(1-L) \varepsilon_t = -(\varepsilon_{t-1} + 1/2 \varepsilon_{t-2} + 1/3 \varepsilon_{t-3} + \dots)$ , it follows that  $T^{-1} \sum \log(1-L) \varepsilon_t \varepsilon_{t-i} = -T^{-1} \sum \varepsilon_{t-i}^2 / i - T^{-1} \sum_{t=i+1}^T \varepsilon_{t-i} \left( \sum_{j=1, j \neq i}^{\infty} \frac{1}{j} \varepsilon_{t-j} \right)$ . It can easily be checked that the later term tends in mean square to zero whereas the LLN for *i.i.d* processes guarantees that the former one converges in probability to  $-\sigma^2/i$ . In turn, this implies that  $\partial \hat{\rho}_e(i) / \partial d \big|_{\lambda=\lambda_0} \xrightarrow{P} -1/i$ , for  $i = 1, \dots, k$ .

The derivatives with respect to the remaining ARMA parameters are given by:

$$\partial \hat{\rho}_e(i) / \partial \phi_j |_{\lambda=\lambda_0} = \frac{1}{T\sigma^2} \sum \frac{-\varepsilon_{t-j}}{\Phi(L)} \varepsilon_{t-i} + o_p(1) \xrightarrow{p} \begin{cases} -\omega_{i-1} & \text{if } j \leq i. \\ 0 & \text{if } j > i. \end{cases}, \quad (42)$$

and:

$$\partial \hat{\rho}_e(i) / \partial \theta_j \xrightarrow{p} \frac{1}{T\sigma^2} \sum \frac{-\varepsilon_{t-j}}{\Theta(L)} \varepsilon_{t-i} + o_p(1) \xrightarrow{p} \begin{cases} -\psi_{i-1} & \text{if } j \leq i \\ 0 & \text{if } j > i. \end{cases} \quad (43)$$

where  $\frac{1}{\Phi(L)} = \sum \omega_i L^i$  and  $\frac{1}{\Theta(L)} = \sum \psi_i L^i$ . Therefore, for finite  $k$ , the Jacobian matrix  $\hat{J}_k(\lambda_0)$  tends to

$$J_k(\lambda_0) = \begin{pmatrix} -1 & 1 & 0 & \dots & 1 & \dots & 0 \\ -1/2 & \omega_1 & 1 & \dots & \psi_1 & \dots & 0 \\ \dots & \dots & \dots & \dots & \dots & \dots & \dots \\ -1/k & \omega_{k-1} & \omega_{k-2} & \dots & \psi_{k-1} & \dots & \psi_{k-q} \end{pmatrix}. \quad (44)$$

Also, since  $\sqrt{T}\hat{\rho}_{ke}(\lambda_0) = \sqrt{T}\hat{\rho}_{k\varepsilon} + o_p(1) \xrightarrow{w} N(0, I_k)$  for fixed  $k$  (Theorem 1), it is straightforward to check that  $\sqrt{T}(\hat{\lambda}_k - \lambda_0)$  is also asymptotically normally distributed with mean equal to zero and variance-covariance matrix given by  $\Xi_k^{-1} = (J'_k(\lambda_0) J_k(\lambda_0))^{-1}$ . The case where  $\mu$  is unknown can be proven along the same lines. ■

In order to prove Theorem 5, where the consistency of the estimator when  $k$  also tends to infinity is established, the sets  $\Lambda_1$  and  $\Lambda_2$  need to be redefined.

**Definition 2** Denote  $\lambda^{(1)} = d$  and  $\Lambda_1^* = \{d : \nabla_1 \leq d \leq \nabla_2\} \times \Lambda^{(-1)}$  if  $\nabla_1 > d_0 - 1/4$  or otherwise  $\Lambda_1^* = \{d : d_0 - 1/4 + \eta \leq d \leq \nabla_2\} \times \Lambda^{(-1)}$  for some  $\eta \in (0, 1/4)$ , if  $\nabla_1 \leq d_0 - 1/4$ , where  $\Lambda^{(-1)}$  is the parameter space of the remaining ARMA parameters.

The reason for introducing this modification is that, since  $k \rightarrow \infty$ , the criterion function contains an increasing number of correlations and the summability of squared correlations of  $FI(d_0)$  processes depends on the value of  $d_0$ . More specifically, whenever  $d_0$  is  $\geq 1/4$ , correlations of  $FI(d_0)$  processes are not square-summable. Following the same reasoning as in the proof of Theorem 3, the integration order of the residuals (obtained after filtering  $y_t$  by  $\Delta^d$  and another ARMA parameters) will be  $(d_0 - d)$ . Then, if  $\nabla_1 > d_0 - 1/4$ , it implies that  $d_0 - d < 1/4$  always, and therefore the associated squared autocorrelations will be summable. With this definition of  $\Lambda_1^*$ , it is ensured that processes whose parameters lie in this set possess squared-summable correlations. The following auxiliary result is needed to prove the theorem.

**Lemma 2** Let  $V_{ke}(\lambda)$  be the criterion function in (12), where  $e_t(\lambda)$  is defined as in (4) or (6) according to the case where the DGP has a known or unknown mean respectively and

let  $\tilde{\lambda} \in \Lambda_1^*$ . Also define  $V_\varepsilon(\tilde{\lambda}) = \sum_{i=1}^{\infty} \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2$ , where  $\rho_{\varepsilon(\tilde{\lambda})}(\cdot)$  are the (population) autocorrelations associated to the non truncated residuals  $\varepsilon(\tilde{\lambda})$ . Let  $k$  be a function of  $T$ , such that  $\lim_{T \rightarrow \infty} k(T) = \infty$  and  $\lim_{T \rightarrow \infty} k(T)/T = 0$ . Then:

1.  $V_\varepsilon(\tilde{\lambda})$  is a continuous function and has a unique minimum at  $\lambda_0$ , such that  $V_\varepsilon(\lambda_0) = 0$ .
2.  $V_{ke}(\lambda)$  is continuous in  $\lambda$ ,  $V_{ke}(\tilde{\lambda})$ ,  $\tilde{\lambda} \in \Lambda_1^*$ , converges in probability to  $V_\varepsilon(\tilde{\lambda})$  and the convergence is uniform.

### Proof of Lemma 2

1. To check the continuity of  $V_\varepsilon(\tilde{\lambda})$  it suffices to show that the partial sum of squared autocorrelations,  $s_n = \sum_{i=1}^n \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2$ , converges uniformly to  $V_\varepsilon(\tilde{\lambda})$  (see Rudin, 1990, Theorem 7.12). Applying the Cauchy criterion, this would be true if and only if for all  $\xi > 0$ , there exists an integer  $N$  such that  $m \geq N$ ,  $n \geq N$  and  $\tilde{\lambda} \in \Lambda_1$ , for which

$$\left| \sum_{i=1}^m \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2 - \sum_{i=1}^n \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2 \right| \leq \xi. \quad (45)$$

Assume, without loss of generality, that  $m > n$ . Then, the left-hand side of (45) is simply  $\sum_{i=n+1}^m \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2$ . Since  $V_\varepsilon(\tilde{\lambda}) < \infty$  for all  $\varepsilon(\tilde{\lambda})$  (recall that  $V_\varepsilon(\tilde{\lambda})$  contains the squared autocorrelations of  $FI$  processes with an order of integration given by  $(d_0 - \tilde{d})$  strictly smaller of  $1/4$ ), there exists an integer  $N$  such that for  $n > N$ ,  $\sum_{i=n+1}^{\infty} \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2 \rightarrow 0$ , which clearly implies that  $\sum_{i=n+1}^m \left( \rho_{\varepsilon(\tilde{\lambda})}(i) \right)^2$  also tends to zero for all  $\tilde{\lambda} \in \Lambda_1^*$ .

It is straightforward to check that it has a unique minimum at  $\lambda_0$ , since  $\varepsilon_t(\lambda)|_{\lambda=\lambda_0} = \varepsilon_t$ , which is an *i.i.d.* process and therefore all its correlations are zero, (which implies that  $V_\varepsilon(\lambda_0) = 0$ ), but presents non-zero autocorrelations for any other value of  $\tilde{\lambda} \neq \lambda_0$  (and therefore  $V_\varepsilon(\tilde{\lambda}) > 0$ ).

2. The continuity property follows along the same lines as in Lemma 1. Let  $V_{ke}(\tilde{\lambda}) = \sum_{i=1}^{k(T)} \left( \hat{\rho}_{\varepsilon(\tilde{\lambda})}(i) \right)^2$ . The convergence in probability is first proven for the autocovariance function and it trivially extends to the autocorrelation one. Following again Proposition 6.3.9 in Brockwell and Davis (1993), it is just needed to show that *i)*  $\sum_{i=1}^j \left( \hat{\gamma}_{\varepsilon(\tilde{\lambda})}(i) \right)^2 \xrightarrow{p} \sum_{i=1}^j \left( \gamma_{\varepsilon(\tilde{\lambda})}(i) \right)^2$  for fixed  $j$ , *ii)*  $\lim_{j \rightarrow \infty} \sum_{i=1}^j \left( \gamma_{\varepsilon(\tilde{\lambda})}(i) \right)^2 =$

$\sum_{i=1}^{\infty} \left( \gamma_{\varepsilon}(\tilde{\lambda})(i) \right)^2$ , and *iii*)

$$\lim_{j \rightarrow \infty} \limsup_{T \rightarrow \infty} P \left( \left| \sum_{i=1}^{k(T)} \left( \hat{\gamma}_{\varepsilon}(\tilde{\lambda})(i) \right)^2 - \sum_{i=1}^j \left( \gamma_{\varepsilon}(\tilde{\lambda})(i) \right)^2 \right| > \epsilon \right) = 0. \quad (46)$$

*i*) follows from Lemma 1; *ii*) follows from noticing that the limit is well defined since autocovariances (and also autocorrelations) of  $FI(d)$  processes are square-summable if  $d < 1/4$ , which is verified by the processes considered here. Therefore  $V_{\varepsilon}(\tilde{\lambda}) < \infty$ . Finally, the probability in (46) is bounded by  $\epsilon^{-1} \left( \sum_{i=j+1}^{k(T)} E \left( \hat{\gamma}_{\varepsilon}(\tilde{\lambda})(i) \right)^2 \right)$  by Markov's inequality. By Proposition 7.3.3. in Brockwell and Davis (1993),

$$\lim_{T \rightarrow \infty} \sum_{i=j+1}^{k(T)} E \left( \hat{\gamma}_{\varepsilon}(\tilde{\lambda})(i) \right)^2 = \sum_{i=j+1}^{\infty} \left( \gamma_{\varepsilon}(\tilde{\lambda})(i) \right)^2,$$

which is finite, since autocovariances are square-summable as stated above. Then, it clearly follows that

$$\lim_{j \rightarrow \infty} \lim_{T \rightarrow \infty} \epsilon^{-1} \sum_{i=j+1}^{k(T)} E \left( \hat{\gamma}_{\varepsilon}(\tilde{\lambda})(i) \right)^2 = 0. \quad (47)$$

which in turn implies that the limit in (46) is also equal zero.

The uniform convergence follows from the pointwise convergence and the same arguments as in Lemma 1.

### Proof of Theorem 5.

The proof of the first part of the theorem is analogous to that of Theorem 3, with Definition 1 and the results of Lemma 1 replaced by Definition 2 and the results of Lemma 2.

*First step.* Define  $\hat{\lambda}_1 = \arg \min_{\lambda \in \Lambda_1^*} V_{ke}(\lambda)$ . Using the same arguments as in Theorem 3, it follows that  $\hat{\lambda}_1 \xrightarrow{P} \lambda_0$ .

*Second step.* Consider the case where  $d_0 > 1/4 + \nabla_1$  and define  $\Lambda_2^* = \{d : \nabla_1 \leq d < d_0 - 1/4 + \eta\} \times \Lambda^{(-1)}$  and  $\hat{\lambda} = \arg \min_{\lambda \in \Lambda} V_{ke}(\lambda)$ . As in the proof of Theorem 3, it is needed to show that for any  $\delta > 0$  the probability,

$$P \left( \left\| \hat{\lambda} - \hat{\lambda}_1 \right\| \geq \delta \right) \leq P \left( \inf_{\lambda \in \Lambda_2} V_{ke}(\hat{\lambda}) - V_{ke}(\hat{\lambda}_1) \leq 0 \right) \quad (48)$$

converges to zero. Using the same arguments as in Theorem 3, it can be checked that  $P \left( V_{ke}(\hat{\lambda}_1) > 0 \right) \rightarrow 0$  and  $P \left( V_{ke}(\lambda) > 0 \right) \rightarrow 1$  for all  $\lambda \in \bar{\Lambda}_2$ , where  $\bar{\Lambda}_2$  is the closure of  $\Lambda_2$  and, since the infimum is contained in this set, it holds that  $P \left( \inf_{\lambda \in \Lambda_2} V_{ke}(\hat{\lambda}) \leq 0 \right) \rightarrow 0$ . The latter arguments imply that the probability in (48) tends to zero.

With respect to the asymptotic distribution of  $\hat{\lambda}$ , application of the mean value theorem to the first-order conditions yields an expression similar to that in (38). In order to apply Theorem 2, it is needed to show that,

$$\left\| \hat{J}_k - J_k \right\| = \left( \sum_{j=1}^{p+q+1} \sum_{i=1}^k \left( \partial \hat{\rho}_e(i) / \partial \lambda_j |_{\lambda=\lambda_0} - J_{ij} \right)^2 \right)^{1/2} = o_p \left( k^{-1/2} \right), \quad (49)$$

where  $J_{ij}$  is the  $-ij$  element of the matrix  $J_k$  defined in (44). Since the first summation is a sum of a finite number of terms, a sufficient condition for (49) is simply,

$$\lim_{T \rightarrow \infty} k^{1/2} \left( \sum_{i=1}^k \left( \partial \hat{\rho}_e(i) / \partial \lambda_j |_{\lambda=\lambda_0} - J_{ij} \right)^2 \right)^{1/2} = 0, \text{ for } j = 1, \dots, p+q+1. \quad (50)$$

To check that this condition is verified, we use again proposition 6.3.9. in Brockwell and Davis (1993). The first two conditions are trivially verified: condition i) follows from results in Theorem 3 while ii) is simply  $\lim_{k \rightarrow \infty} 0$ , which is obviously equal to zero. The last condition can be shown along the same lines as Lemma 2, (2), noticing that  $TE \left( \left( \partial \hat{\rho}_e(i) / \partial \lambda_j |_{\lambda=\lambda_0} - J_{ij} \right)^2 \right) = O_p(1)$ , and that the corresponding expectations are summable in  $i$  (for instance, the limit of the expectation for the derivative with respect to  $d$  is equal to  $\mu_4/i^2$ , where  $\mu_4 = E(\varepsilon^4)$ , and therefore, summable over  $i$ ). Finally, the result follows just by noticing that  $k/T \rightarrow 0$ . By Theorem 2 and taking into account the consistency of  $\hat{\lambda}$ , it follows that

$$\sqrt{T} \left( \hat{\lambda} - \lambda_0 \right) \xrightarrow{w} N \left( 0, \Xi^{-1} \right)$$

where  $\Xi = \lim_{k \rightarrow \infty} \Xi_k$ . It just remains to be shown that the matrix  $\Xi$  is in fact the Fisher information matrix. Recall from the proof of Theorem 4 that the limit of the Jacobian matrix was given by

$$J_k(\lambda_0) = \begin{pmatrix} -1 & 1 & 0 & \dots & 1 & \dots & 0 \\ -1/2 & \omega_1 & 1 & \dots & \psi_1 & \dots & 0 \\ \dots & \dots & \dots & \dots & \dots & \dots & \dots \\ -1/k & \omega_{k-1} & \omega_{k-2} & \dots & \psi_{k-1} & \dots & \psi_{k-q} \end{pmatrix}. \quad (51)$$

Given the previous result, it is easy to obtain that

$$\Xi = \begin{pmatrix} \pi^2/6 & \Pi' \\ \Pi & \Xi_{pq} \end{pmatrix}$$

where  $\Pi = (\pi_\omega(0), \dots, \pi_\omega(p-1), \pi_\psi(0), \dots, \pi_\psi(q-1))$  and

$$\pi_\omega(j) = \sum_{i=0}^{\infty} \frac{\omega_i}{j+i+1}, \quad \pi_\psi(j) = \sum_{i=0}^{\infty} \frac{\psi_i}{j+i+1}.$$

$\Xi_{pq}$  is the Fisher information matrix corresponding to pure ARMA processes. Then,  $\Xi$  coincides with the Fisher information matrix for ARFIMA processes (see for instance Li and McLeod, 1986 or Fox and Taqqu 1996).■

**Proof of Theorem 6**

Let  $\hat{\rho}_{ke}(\hat{\lambda})$  be the vector defined in (19). A first-order Taylor's series expansion of  $\hat{\rho}_{ke}(\hat{\lambda})$  around  $\lambda_0$  yields:

$$\hat{\rho}_{ke}(\hat{\lambda}) = \hat{\rho}_{ke}(\lambda_0) + \frac{\partial \hat{\rho}_{ke}(\lambda^*)}{\partial \lambda'} (\hat{\lambda} - \lambda_0) \quad (52)$$

By Theorem 1,  $\sqrt{T}\hat{\rho}_{ke}(\lambda_0) = \sqrt{T}\hat{\rho}_{ke} + o_p(1) \xrightarrow{w} N(0, I_k)$  for  $d > -0.75$ . The joint distribution of  $\left(\hat{J}_k(\lambda_0) (\hat{\lambda} - \lambda_0), \hat{\rho}_{ke}\right)'$  is given by:

$$\sqrt{T} \begin{pmatrix} \hat{J}_k(\lambda_0) (\hat{\lambda} - \lambda_0) \\ \hat{\rho}_{ke} \end{pmatrix} = \begin{pmatrix} -\hat{J}_k(\lambda_0) \left(\hat{J}'_k(\lambda_0) \hat{J}_k(\lambda_0)\right)^{-1} \hat{J}'_k(\lambda_0) \\ I_k \end{pmatrix} \sqrt{T}\hat{\rho}_{ke} + o_p(1)$$

$$\xrightarrow{w} N(0, \Upsilon),$$

where:

$$\Upsilon = \begin{pmatrix} -J_k(\lambda_0) (J'_k(\lambda_0) J_k(\lambda_0))^{-1} J'_{\lambda_0 k} & J_k(\lambda_0) (J'_k(\lambda_0) J_k(\lambda_0))^{-1} J'_k(\lambda_0) \\ J_k(\lambda_0) (J'_k(\lambda_0) J_k(\lambda_0))^{-1} J'_k(\lambda_0) & I_k \end{pmatrix}.$$

Since joint normality holds, any linear combination of  $\left(\hat{J}_k(\lambda_0) (\hat{\lambda} - \lambda_0), \hat{\rho}_{ke}\right)$  will also be normal. Taking into account expression (52) it follows that:

$$\sqrt{T}\hat{\rho}_{ke}(\hat{\lambda}) \xrightarrow{w} N\left(0, I_k - J_k(\lambda_0) (J'_k(\lambda_0) J_k(\lambda_0))^{-1} J'_k(\lambda_0)\right). \blacksquare$$

## APPENDIX 2

This appendix reports the results of some Monte-Carlo experiments not included in the main text. Table 12 presents the mean and standard deviation of correlations at different lags associated to the (truncated) residual process  $e_t(d_0)$  for different values of  $d_0$ . To obtain these residuals, the following procedure has been implemented: processes of the form  $\Delta^{\varphi_0} y_t = \varepsilon_t$  were generated for a large sample size equal to  $n+T$  and then the first  $n = 1000$  observations were rejected. The last  $T$  observations were integrated an integer number of times,  $m_0$ . Truncated residuals were computed by applying the finite filter  $\sum_{i=0}^{t-1} \pi_i L^i$  to the process  $y_t$ , where the coefficients  $\{\pi_i\}$  come from the expansion in powers of  $L$  of the polynomial  $(1-L)^{d_0}$ . Again the results in finite samples confirm the asymptotic results. It can be seen that for invertible processes ( $d_0 > -1$ ), estimated correlations behave correctly in the sense that, as expected, zero mean and unit variance is found. Nevertheless, for non-invertible processes, residuals correlations do not provide consistent estimates of the innovation correlations.

**Table 12.** Residual Autocorrelations

$d_0$	-1.2	-1.0	-0.7	0.4	0.8	1.4
$T = 100$						
mean( $\sqrt{T}\hat{\rho}_e(1)$ )	0.801	-0.136	-0.063	-0.061	-0.128	-0.068
std( $\sqrt{T}\hat{\rho}_e(1)$ )	1.435	0.993	0.964	0.966	0.955	0.979
mean( $\sqrt{T}\hat{\rho}_e(2)$ )	0.804	-0.085	-0.022	-0.037	-0.080	-0.027
std( $\sqrt{T}\hat{\rho}_e(2)$ )	1.386	0.971	0.977	0.979	0.995	0.990
mean( $\sqrt{T}\hat{\rho}_e(5)$ )	0.793	-0.046	0.009	-0.006	-0.042	-0.002
std( $\sqrt{T}\hat{\rho}_e(5)$ )	1.361	0.956	0.966	0.964	0.971	0.974
$T = 400$						
mean( $\sqrt{T}\hat{\rho}_e(1)$ )	3.020	-0.006	0.010	0.047	-0.001	0.042
std( $\sqrt{T}\hat{\rho}_e(1)$ )	3.248	0.979	1.001	0.982	0.979	0.983
mean( $\sqrt{T}\hat{\rho}_e(2)$ )	2.926	-0.045	-0.031	-0.005	-0.039	-0.005
std( $\sqrt{T}\hat{\rho}_e(2)$ )	3.170	1.013	1.027	1.016	1.013	1.019
mean( $\sqrt{T}\hat{\rho}_e(5)$ )	2.915	-0.006	0.002	0.026	-0.002	0.029
std( $\sqrt{T}\hat{\rho}_e(5)$ )	3.149	0.968	1.031	0.958	0.969	0.964