WHICH MODEL TO MATCH?

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Abstract

The asymptotic efficiency of indirect estimation methods, such as the efficient method of moments and indirect inference, depends on the choice of the auxiliary model. To date, this choice has been somewhat ad hoc and based on an educated guess. In this article we introduce a class of information criteria that helps the user to optimize the choice between nested and non–nested auxiliary models. They are the indirect analogues of the widely used Akaike–type criteria. A thorough Monte Carlo study based on two simple and illustrative models shows the usefulness of the criteria.

Keywords: Indirect inference, efficient method of moments, auxiliary model, information criteria, asymptotic efficiency.

JEL classification: C13, C52.
Resumen

La eficiencia asintótica de los estimadores de inferencia indirecta, tales como el método eficiente de los momentos, depende de la elección del modelo auxiliar. Hasta ahora, esta elección era ad hoc y basada en criterios subjetivos. En este artículo introducimos una clase de criterios de información que ayuda al usuario a escoger entre modelos anidados y no anidados. Esta clase es la análoga a los ampliamente utilizados criterios del tipo Akaike. Un detallado estudio de Monte Carlo basado en dos modelos simples e ilustrativos muestran la utilidad de los criterios.

Palabras clave: inferencia indirecta, método eficiente de los momentos, modelo auxiliar, criterios de información, eficiencia asintótica.

Códigos JEL: C13, C52.
1 Introduction

The title of this article intentionally resembles to Gallant and Tauchen (1996)’s *Which moments to match?* They, Gouriéroux et al. (1993), and Smith (1993) introduce an inference method based on matching sample and theoretical features of an auxiliary model in order to estimate the parameters of the model of interest. Gouriéroux et al. (1993) call it Indirect Inference, Smith (1993) Simulated Quasi Maximum Likelihood, and Gallant and Tauchen (1996) the Efficient Method of Moments (EMM). We will denote these methods generically as *indirect methods*.

Because of their flexibility, applications of indirect methods in all areas of economics abound.\(^1\) Provided that the auxiliary estimators converge to some point in their space and that the so–called binding function is injective, then the general theory shows that any auxiliary model is valid. Validity in this context means that the estimated parameters are, under fairly weak regularity conditions, consistent, asymptotically Gaussian, and with a variance–covariance matrix that is a function of the binding function and the Fisher information matrix of the auxiliary model.

In practice, a good auxiliary model is a model that, in some sense, is close to the one of interest. Gallant and Tauchen (1996) introduce the SNP (Semi Non–Parametric) score generator as universal auxiliary model. It is based on Hermite expansions that can, in principle, span the score vector of any model of interest. The SNP score generator is a class of auxiliary models since there are as many models as terms in the expansion; so a choice has to be made. An alternative to SNP is to use an auxiliary model that is much bigger (and hence with many more parameters) than the model of interest, on the grounds that the bigger the auxiliary model, the better the approximation to the model of interest. In the limit, this leads to the sieve–based approach of Nickl and Potscher (2010) and Blasques (2011). Though theoretically it is a reasonable argument, in practice we face limited information given by the sample size, meaning that the finite sample properties of the estimated parameters may deteriorate sensibly with the dimension of the auxiliary model. This is the classical trade–off between model fit and the number of parameters.

To the best of our knowledge no systematic analysis has been done on how to choose the auxiliary model within the general theory of indirect methods. In this article we propose information criteria that can be used for choosing between a set of nested and non–nested auxiliary models. They are the indirect analogues of the widely used Akaike–type model information criteria (among those, the most widely used are AIC of Akaike (1974), BIC of Schwarz (1978), and HQIC of Hannan and Quinn (1979)). The indirect auxiliary model selection criteria (that we denote with subscript $IM$ for Indirect Methods) are based on the likelihood of the estimated parameters of the model of interest, which is centered

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\(^1\)A non–exhaustive list is: discrete stochastic volatility models (Monfardini (1998), Gallant et al. (1997), Lombardi and Calzolari (2009), and Corsi and Reno (2012)), continuous stochastic volatility models (Chernov et al. (2003) and Grassi and de Magistris (2012)), diffusion models (Engle and Lee (1996), Broze et al. (1998)), threshold models (Guay and Scaillet (2003)), heavy tailed distributions (Lombardi and Veredas (2009), Garcia et al. (2011)), DSGE models (Dridi et al. (2007)), survival data models (Jiang and Turnbull (2003)), stochastic agent–based models (Jang and Lux (2012)), labor income risk (Guvenen and Smith (2010)), discrete choice models (Keane and Smith (2003)), heteroskedastic factor models (Calzolari et al. (2008)), learning models (Czellar and Calvet (2011)), and price contingent claims in derivative markets (Phillips and Yu (2009)).
Gaussian regardless of the choice of the auxiliary model. The difference therefore comes in terms of precision, i.e. the variance–covariance matrix. But to account for the trade–off between the fit of the auxiliary model to the model of interest and its dimension, we need a penalty that depends on the number of parameters of the auxiliary model.

Our work relates to the existing literature on Generalized Method of Moments (GMM) for choosing the moment conditions (MC). Two issues arise in the GMM context: the choice of MC that ensures consistency of the estimates, and the choice of no redundant moment conditions. The former is related to weak identification, or to the case where there is insufficient information to yield a consistent estimator. Andrews (1999) proposes GMM analogues of the AIC, BIC and HQIC (which we denote with subscript $GMM$) for consistently selecting the correct MC.\(^2\) As for the choice of no redundant MC, it relates with moment conditions that provide no incremental information about the parameters of the model of interest. Hall et al. (2007) introduce the Relevant Moment Selection Criteria (RMSC) where the word relevant denotes the subset of MC that are asymptotically efficient but that contain no redundant moment conditions. Though redundant moment conditions have no asymptotic impact, Hall and Peixe (2003) report simulation evidence that their inclusion can lead to a poor approximation of the limiting distribution to the finite sample behavior.

In the framework of indirect methods, the choice of the auxiliary model presents no concerns in terms of consistency, provided the conditions stated above. The closest to our work is therefore Hall et al. (2007). They propose RMSC based on the observation that the entropy of the limiting distribution of the GMM estimator provides a metric for information of the moment conditions. These ideas were also used by Hall et al. (2011) in the context of DSGE models, where the auxiliary model are impulse response functions (IRF hereafter) steaming from a VAR. This is alike to Dridi et al. (2007) that estimate a DSGE model with sequential partial indirect inference.

The indirect inference criteria that we present can also be used for choosing between and within other matching inference methods, such as GMM, the Method of Simulated Moments (MSM; McFadden (1989), Pakes and Pollard (1989), and Duffie and Singleton (1993)) or the Method of Simulated Quantiles (MSQ; Dominicy and Veredas (2012)). As long as the functions to match lead to consistent estimators, the only difference between them is in terms of the variance–covariance matrix, as in the indirect methods. In other words, the information criteria only depend on the variance–covariance matrix of the estimators and the number of functions to match.

We proceed with a thorough Monte Carlo study on two simple, yet important and illustrative, models: a moving average (MA) of order one and a stochastic volatility (SV) model. In this context of indirect methods, the former has been studied by Gouriéroux et al. (1993), Ghysels et al. (2003), and Genton and Ronchetti (2003) while the latter by Monfardini (1998) and Calzolari et al. (2004). For the MA, we choose as auxiliary models an AR($r$) and the true model. The purpose of this study is to investigate the performance of the criteria when choosing among the true model and nested models. The

\(^2\)See also Andrews and Lu (2001), Kitamura (2002), Hong et al. (2003) and Hong and Preston (2005) for related work on moment selection criteria, and Konishi and Kitagawa (1996) for a generalized information criteria.
criteria choose the true model as best auxiliary model. The second best is the AR model with lag order that increases with the magnitude of the moving average parameter. As for the SV model, we consider two auxiliary models: ARCH(r) and GARCH(1,1) on the observations, and AR(r) and ARMA(1,1) on the log of the square of the observations. The purpose of this study is to investigate the performance of the criteria when choosing among nested and non–nested auxiliary models. The criteria choose r that increases with the SV autoregressive parameter, and always prefer the ARCH to the AR model.

The remaining sections are laid out as follows. Section 2 introduces the notation and the basic background of indirect methods, while Section 3 describes the criteria. Section 4 covers the Monte Carlo study. Conclusions are touched upon in Section 5. Assumptions, technical lemmas, proofs, and tables are relegated to the appendices.

2 Indirect methods

The model of interest We consider a \( n \times 1 \) random vector \( Y \) that follows a model \( M(y, \theta) \), where \( \theta \) denotes the \( p \)-dimensional vector of unknown parameters that are in an interior point of the parameter set \( \Theta \subset \mathbb{R}^p \), and \( y \in \mathcal{Y} \subset \mathbb{R}^n \) is a realization of \( Y \). The model is defined by a parametric probability law

\[
M(y, \theta) = \{ P_{\theta}(y), y \in \mathcal{Y} \subset \mathbb{R}^n, \theta \in \Theta \subset \mathbb{R}^p \}. \tag{1}
\]

The true value of the parameters is denoted by \( \theta_0 \) and it is an interior point of \( \Theta \). We define the Fisher information matrix for this model as \( I(\theta_0) \). We assume to have \( T \) realizations of \( Y \) and we denote as \( y_t = (y_{1t} \ldots y_{nt})' \) the \( n \times 1 \) vector of a generic realization with \( t = 1, \ldots, T \).

The class of models we consider is broad, covering those quoted in the introduction, and including static and dynamic models, with unobserved components, trends, and various innovation terms. Each process and model has its own assumptions. Establishing general hypothesis on \( Y \) and \( M(y, \theta) \) is therefore idle, beyond A.1 in Appendix A on \( \theta \) and its space. It is however important that the model is such that, for a given value of the parameter vector, simulated draws of \( y_1(\theta), \ldots, y_T(\theta) \) can be obtained. This typically implies that \( M(y, \theta) \) contains at least one i.i.d. innovation with known distribution.

The auxiliary model Consider the auxiliary model for \( Y \)

\[
M^a(y, \beta) = \{ P_{\beta}(y), y \in \mathcal{Y} \subset \mathbb{R}^n, \beta \in \mathcal{B} \subset \mathbb{R}^q \}
\]

that depends on a \( q \)-dimensional vector of parameters \( \beta \). We assume that for model \( M^a(y, \beta) \) the log–likelihood \( \ell^a(\beta; y_t) \) is known and available analytically. We define \( \hat{\beta}_T \) as the Quasi Maximum Likelihood (QML) estimator of the true value of the parameters \( \beta_0 \). Under the standard regularity conditions of QML estimation (see also assumptions A.1–A.4 in Appendix A), \( \hat{\beta}_T \) has the following limiting distribution

\[
\sqrt{T}(\hat{\beta}_T - \beta_0) \sim \mathcal{N}(0_q, (J(\beta_0)I^{-1}(\beta_0)J(\beta_0))^{-1}), \quad \text{as } T \to \infty,
\]

where \( 0_q \) is a \( q \times 1 \) vector of zeros, \( I(\beta_0) \) is the \( q \times q \) Fisher information matrix of the auxiliary model, and \( J(\beta_0) \) is minus the expectation of the Hessian of the log–likelihood.
For a given value of \( \theta \), we simulate \( H \) paths of length \( T \): \( y_t^h(\theta), \ldots, y_T^h(\theta), h = 1, \ldots, H \), from the model of interest \( M(y, \theta) \) and we estimate

\[
\hat{\beta}_{HT}(\theta) = \arg \max_{\beta \in \mathcal{B}} \frac{1}{HT} \sum_{h=1}^{H} \sum_{t=1}^{T} c^*(\beta, y_t^h(\theta)).
\]  

(2)

Under assumptions A.1–A.4 in Appendix A, this is a consistent and asymptotically Gaussian estimator (as \( T \to \infty \)) of the so-called binding function \( b(\theta) \) linking \( \theta \) to \( \beta \) and such that \( b(\theta_0) = \beta_0 \). Finally, notice that we could simulate paths of length \( T^* \neq T \) (typically \( T^* > T \)) and therefore both the estimated auxiliary parameters and their estimated asymptotic variance–covariance matrix will be functions of \( T^* \) instead of \( T \). For simplicity of notation we omit this distinction in what follows.

**Indirect estimation** The indirect estimator is defined as

\[
\hat{\theta}_{HT,IM,a,q} = \arg \min_{\theta \in \Theta} \left( \hat{\beta}_{HT}(\theta) - \beta_T \right)' \Omega \left( \hat{\beta}_{HT}(\theta) - \beta_T \right),
\]  

(3)

where \( \Omega \) is a \( q \times q \) weighting matrix defining the metric. We make explicit the dependence of the estimated parameter \( \hat{\theta}_{HT,IM,a,q} \) on the auxiliary model \( a \) and the dimension of its parameter vector \( q \). The \( \beta \)--based optimization (3) was introduced by Gouriéroux et al. (1993). Alternatively we may consider a \( h(\beta) \)--based optimization, for \( h \) an injective and continuously differentiable function with respect to \( \beta \), and after a suitable re–parametrization of the weighting matrix by means of the delta method. The vector \( h(\beta) \) is of dimension \( q' \) that can be larger or equal to \( q \). In the score–based optimization of Gallant and Tauchen (1996) \( q' = q \), while in the IRF–based optimization of Hall et al. (2011) \( q' > q \). A special case is when the auxiliary model is estimated under constraints on the parameter space, i.e. constrained QML that imply the estimation of multipliers (Calzolari et al. (2004)). The vector \( \beta \) is expanded to dimension \( q' > q \) as to include the multipliers. But it is \( q \) that enters in the penalty. If the constraints are not binding, the multipliers are zero, and inference is as unconstrained. If, by contrast, the constraints bind, the bound parameters contain no information about \( \theta \), as it goes through the associated multipliers. In the sequel, we consider the case of the \( \beta \)--based optimization, and hence the penalty of the criteria is a function of \( q \). If, instead, the \( h(\beta) \)--based optimization is used, the criteria are easily adapted by replacing \( q \) by \( q' \).

Under assumptions A.1–A.7 in Appendix A and the appropriate model assumptions, the estimators are consistent with limiting distribution

\[
\sqrt{T}(\hat{\theta}_{HT,IM,a,q} - \theta_0) \sim N(0, W_{HT, IM, a,q}(\theta_0)), \quad \text{as} \quad T \to \infty.
\]

The asymptotic variance–covariance matrix depends on the true value of the parameters, the number of simulated paths, and the weighting matrix. The optimal choice for the latter is \( \Omega^* = J(\beta)I^{-1}(\beta)J(\beta) \) so that

\[
W_{HT, IM, a,q}(\theta) = \left(1 + \frac{1}{H} \right) \left[ \frac{\partial b(\theta)}{\partial \theta} J(\beta) I^{-1}(\beta) J(\beta) \frac{\partial b(\theta)}{\partial \theta'} \right]^{-1}.
\]

Since the value of \( H \) is arbitrarily chosen, we consider \( H \to \infty \) and, as in the next section we need further notation, hereafter we skip the subindexes \( H \) and \( \Omega^* \), so \( \hat{\theta}_{HT,IM,a,q} \).
becomes $\hat{\theta}_{T,IM,a,q}$ and $W_{H,\Omega^*,a,q}(\theta)$ becomes

$$W_{a,q}(\theta) = \left[ \frac{\partial b(\theta)}{\partial \theta} J(\beta) \Gamma^{-1}(\beta) J(\beta) \frac{\partial b(\theta)}{\partial \theta'} \right]^{-1}.$$  

(4)

A plethora of theoretical developments and improvements of the original methods have been proposed. What follows is a selection of an extensive and growing literature. Dhaene et al. (1998) introduce a notion of indirect encompassing and non-nested hypotheses tests using indirect methods. That is, the case where not only the relevant binding function does not have closed form, but also at least one of the competing models. Billio and Monfort (2003) propose kernel-based conditional expectations for the binding functions, which have the advantage that no optimization step is involved in the computation of the binding function. They also propose two criteria specific for conditional expectations as binding functions. Another area of research within indirect methods concerns robustification. Genton and Ronchetti (2003) develop robust indirect inference by deriving the influence function of the indirect estimator, as well as the level and power influence function of indirect tests. In a similar spirit, Trojani and Ortelli (2005) focus on robust EMM estimation of a general parametric time series stationary processes. Czellar and Ronchetti (2010) propose robust over-identification tests that exhibit better finite sample accuracy than classical tests. Partial and semi-nonparametric indirect methods have also been introduced. Dridi et al. (2007) introduce the sequential partial indirect inference approach, in the sense that the model of interest is partially well specified. They show the usefulness of the sequential method on the calibration of dynamic stochastic general equilibrium (DGSE) models. Also motivated by DSGE models, Blasques (2011) presents semi-nonparametric indirect methods, i.e. a sieve class of estimator, on the grounds that, at population level, an infinite dimensional auxiliary model may be needed for the indirect estimation of many economic models. Indirect methods have also been used for bias correction (see Gouriéroux et al. (2000) for general time time series models, Ghysels et al. (2003) for a MA(1), and Gourieroux et al. (2010) for dynamic panel data models). The advantage with other bias correction techniques is that indirect methods do not require a given explicit form for the bias function of its expansion, as it is calibrated via simulations. Last, Phillips (2012) provides new limit theory for the delta method and the continuous mapping theorem for cases when the problem involves sample functions that depend on the sample size, or when the quantity of interest appears in an implicit functional form, as it happens in indirect methods.

3 The indirect information criteria

The principle Consider two auxiliary models $\mathcal{M}^{a_1}(y, \beta_1)$ and $\mathcal{M}^{a_2}(y, \beta_2)$ with $\beta_1$ and $\beta_2$ of dimension $q_1$ and $q_2$ respectively. Under the optimal choice of the weighting matrices and as $T, H \to \infty$ we have

$$\sqrt{T}(\hat{\theta}_{T,IM,a,q_1} - \theta_0) \sim N(0, W_{a_1,q_1}(\theta_0)) \quad \text{and} \quad \sqrt{T}(\hat{\theta}_{T,IM,a,q_2} - \theta_0) \sim N(0, W_{a_2,q_2}(\theta_0)).$$
Both estimators are consistent and the limiting distribution of $\sqrt{T}(\hat{\theta}_{T,IM,a,q} - \theta_0)$ is centered Gaussian. The difference between them comes in terms of precision, i.e. the variance–covariance matrix. Choices of $M^a$ and $q$ cover both nested and non–nested auxiliary models. For simplicity of notation we use the index $a$ to indicate a generic auxiliary model $M^a$. In particular, if the models are nested $M^{a_1} \subset M^{a_2}$ (in the sense that they belong to the same auxiliary class but $q_1 < q_2$), we use the same index $a = a_1 = a_2$ for both. If the models are non–nested, there is no well defined inclusion relation between $M^{a_1}$ and $M^{a_2}$, $q_1$ and $q_2$ may or may not be equal, and we say $a_1 \neq a_2$. Finally, we define $A_{q_{\text{max}}}$ as the set of all possible auxiliary models satisfying assumptions A.1–A.7 in Appendix A, where $q_{\text{max}}$ is the largest dimension of the auxiliary parameter we allow for. Each element of $A_{q_{\text{max}}}$ is made of a couple $(a, q)$ indicating the chosen auxiliary model and the dimension of the corresponding auxiliary parameter.

We say that a matrix $A$ is larger than another $B$, and we denote it as $A \succ B$, when $||A|| > ||B||$, where $|| \cdot ||$ is a norm (such as the Frobenius norm $||A||^2 = \text{Tr}(AA')$).

Consider two nested (thus belonging to the same class $1$) auxiliary models with $q_1 < q_2$. Parzen (1959) and Carrasco and Florens (2000) show that

$$W_{a,q_1}(\theta_0) \succ W_{a,q_2}(\theta_0) \succ W_{a,q_{\text{max}}}(\theta_0).$$

That is, the more auxiliary parameters, the more information and the smaller is the asymptotic variance–covariance of the estimated parameters. Whatever the value of $q_{\text{max}}$, the lower bound for the auxiliary models is always larger than the Cramér–Rao lower bound, i.e. $W_{a,q_{\text{max}}}(\theta_0) \succ I^{-1}(\theta_0)$ for any $a$ and $q_{\text{max}}$ (Gouriéroux and Monfort (1996)). Hence, the optimal choice is the largest possible auxiliary model. But this is not the case when we replace $\theta_0$ by $\hat{\theta}_{T,IM,a_1,q_1}$ as we have to take into account the increasing estimation error due to the increasing number of estimated parameters. Some penalization that is a function of $q$ is thus required.

**Simulating the estimated parameters** For computing the indirect criteria that we propose below, we need realizations of the random vector $\sqrt{T}(\hat{\theta}_{T,IM,a,q} - \theta_0)$ that we do not have since $\theta_0$ is not observed. But they are straightforward to simulate. Let $X_{a,q} \equiv \sqrt{T}(\hat{\theta}_{T,IM,a,q} - \theta_0)$, so that $X_{a,q} \sim N(0, W_{a,q}(\theta_0))$, as $T \to \infty$. Let $\hat{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q})$ be the estimator of $W_{a,q}(\theta_0)$ (Gouriéroux et al. (1993)). We know that

$$p \lim_{T \to \infty} \hat{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q}) = W_{a,q}(\theta_0).$$

We can therefore simulate a sample of length $N$ from a $p$–dimensional centered Gaussian distribution with variance–covariance matrix $\hat{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q})$. Let us denote the pdf of $X_{a,q}$ by $f(x_{a,q}, \hat{\theta}_{T,IM,a,q}))$, and the $i$–th simulated observation of $X_{a,q}$ as the $p$–dimensional vector $x_{a,q,i} = (x_{1,a,q,i} \ldots x_{p,a,q,i})'$, with $i = 1, \ldots, N$.

Notice that $n$, $T$ and $N$ are three independent quantities with different meanings. The first one, $n$, is the dimension of the random vector $Y$, while $T$ is the sample length of the

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3For the sake of simplicity, in this section to compare two classes of auxiliary models and for each class we consider the same maximum number of parameters $q_{1,\text{max}} = q_{2,\text{max}} = q_{\text{max}}$. It is however possible that $q_{1,\text{max}} \neq q_{2,\text{max}}$, as in the SV model in Section 4.

4Alternatively, A is larger than a matrix B if A – B is a positive definite matrix.
observed data. Once the data is given, \( n \) and \( T \) are fixed. By contrast, \( N \) is the sample length of the simulated vector \( X_{a,q} \), and can be chosen large enough for the Gaussian limiting distribution to hold. If we define the sample variance–covariance matrix of \( X_{a,q} \) as

\[
\tilde{W}_{N,a,q} = \frac{1}{N} \sum_{i=1}^{N} x_{a,q,i} x_{a,q,i}^\prime,
\]

then the following holds

\[
p \lim_{T \to \infty} \left( p \lim_{N \to \infty} \tilde{W}_{N,a,q} \right) = p \lim_{T \to \infty} \left( \tilde{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q}) \right) = W_{a,q}(\theta_0).
\]

### The criterion

The indirect Akaike information criterion, \( AIC_{IM}(a,q) \), is given in the following proposition.

**Proposition 1** Given the models of Section 2 and the assumptions A.1–A.8 and A.10 in Appendix A, the optimal auxiliary model is the one within the model class \( \tilde{a} \) and auxiliary parameter dimension \( \tilde{q} \) such that

\[
(\tilde{a}, \tilde{q}) = \arg \min_{(q,a) \in A_{\text{max}}} AIC_{IM}(a,q),
\]

where

\[
AIC_{IM}(a,q) = -2 \sum_{i=1}^{N} \log \phi(x_{a,q,i} ; \tilde{W}_{N,a,q}) + \frac{q(q+1)}{2}.
\]  

(5)

**Proof:** See Appendix B.

Three remarks are in order.

First, we give the intuition for the nested case, i.e. when the class of auxiliary models, \( a \), is fixed. Figure 1 gives a geometric interpretation of this problem. The diagram is divided in two rows and three columns. The rows correspond to the largest dimension of the auxiliary model (top row) and a given dimension (bottom). The columns correspond to the variance–covariance matrices in population (right), and the sample and simulated counterparts (middle and left respectively). The calligraphic letters denote vectors.\(^5\) Asymptotically the auxiliary model that is the closest to the model of interest has the maximum possible dimension with variance–covariance matrix of the estimated parameters equal to \( W_{a,q_{\text{max}}}(\theta_0) \), with \( q_{\text{max}} \) large enough. In practice however, \( W_{a,q_{\text{max}}}(\theta_0) \) is unknown and the variance–covariance matrix is estimated for a given \( q < q_{\text{max}} \). \( \tilde{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q}) \). The aim is to find the value of \( q \) that minimizes the distance between \( W_{a,q_{\text{max}}}(\theta_0) \) and \( \tilde{W}_{T,a,q}(\hat{\theta}_{T,IM,a,q}) \). This value is not necessarily \( q_{\text{max}} \), for fixed \( T \), as when \( q \) increases the number of parameters to estimate increases and the estimation error increases accordingly. However, since we use simulated samples \( X_{a,q} \), we need to use asymptotic results for large values of the sample size \( N \). For this reason we have to consider the sample covariance \( \tilde{W}_{N,a,q} \), and thus the left column in the diagram.

\(^5\)The right and middle columns of this diagram are also the base of \( AIC_{ML} \) and \( AIC_{GMM} \).
The diagram is divided in two rows and three columns. The rows correspond to the largest dimension of the auxiliary model (top row) and a given dimension (bottom). The columns correspond to the variance–covariance matrices in population (right), and the sample and simulated counterparts (middle and left respectively). The calligraphic letters denote vectors.

What follows is a sketch of the proof for the nested case. Let \( w = \text{vech}(W) \) and \( Q = (w_{0,a,q_{\text{max}}} - \hat{w}_{T,a,q}) \). Then the distance between \( W_{0,a,q_{\text{max}}} \) and \( \hat{W}_{T,a,q} \) can be written as the norm squared

\[
Q^2 = ||w_{0,a,q_{\text{max}}} - \hat{w}_{T,a,q}||_J^2,
\]

where the norm is based on Hessians \( J \) as defined in Appendix B. We want the auxiliary model \((a,q)\) that minimizes the distance \( Q^2 \). From the geometry of Figure 1, this distance equals

\[
Q^2 = (P - G)^2 = P^2 + G^2 - 2GP.
\]

From Lemma 1 in Appendix B, the term \( G^2 \) is distributed as \( \chi^2_{q(q+1)/2} \), while the last term has expectation zero for large \( N \). Finally, by Pythagora’s theorem we have

\[
P^2 = (A + C)^2 + B^2.
\]

Now \( A + C \) does not depend on \( q \) so it does not play any role in the criterion and we can treat it as a constant, say \( k \), and \( B^2 \) gives the log–likelihood term computed in \( \hat{W}_{N,a,q} \). Therefore, by letting \( N \) diverging and taking expectations of \( (7) \), we can write

\[
\]

\[
\simeq k + \frac{2}{N} \sum_{i=1}^{N} \log \frac{\phi(x_{i,a,q_{\text{max}}}, \hat{W}_{N,a,q_{\text{max}}})}{\phi(x_{i,a,q}, \hat{W}_{N,a,q})} + \frac{q(q+1)}{2N}.
\]

By multiplying the last expression by \( N \) and keeping only the terms that depend on \((a,q)\) we have the criterion. Summing up, \( Q^2 \) depends on i) the distance between \( W_{a,q_{\text{max}}}^{\theta_0} \) and its projection in a \( q \) dimensional subspace \( W_{a,q}^{\theta_0} \), i.e. the term \( B^2 \), and ii) the distance between \( \hat{W}_{N,a,q} \) and the estimated variance–covariance matrix of the auxiliary model \( \hat{W}_{T,a,q}^{\theta_{T,IM,a,q}} \), i.e. the term \( G^2 \). While the first distance decreases as \( q \) increases, the second distance is an estimation error –or a penalty– that increases with \( q \).

While this is the sketched proof for the nested case, \( AIC_{IM} \) compares nested and non–nested models, which implies that the best model is the true model, i.e. the one with
asymptotic variance covariance matrix $I^{-1}(\theta_0)$. But since this is never achievable, the geometric argument cannot be made. For this reason a more involving proof than original Akaike’s is needed –as it is shown in detail in Appendix B– and that is valid for the nested and non–nested cases.

Second, there are important differences between $AIC_{LM}$ and the $ML$ and $GMM$ counterparts. First, the likelihood term in $AIC_{LM}$ is of the estimated parameters of the model of interest, which is centered Gaussian regardless of the choice of the auxiliary model and the model of interest (under assumptions A.1–A.8 in Appendix A and the appropriate model assumptions). By contrast, in $AIC_{ML}$ the likelihood term depends on the model of interest and in $AIC_{GMM}$ depends on the moment conditions. Second, in $AIC_{LM}$ the penalty term depends on the parameters of the auxiliary model, while in $AIC_{ML}$ it depends on the parameters of the model of interest. Third, the dimension of the parameters of the model of interest is constant over choices of the auxiliary model, whilst in $AIC_{ML}$ the dimension varies with the model.

Finally, as it happens with $AIC_{ML}$ and $AIC_{GMM}$, $AIC_{LM}$ is not consistent when $N \to \infty$, in the sense that it tends to select large models because the penalty does not depend on the number of observations. Indeed when $N \to \infty$ the first term on the right hand side of (5) diverges to $-\infty$, thus making the penalty term not effective. As a consequence, $AIC_{LM}$ will not necessarily select the true model if included in the set of all possible auxiliary models, and in this sense we say that it is not consistent. By modifying the penalty, we can generalize $AIC_{LM}$ to obtain consistent criteria.

We define as $(a^*, p)$ the couple corresponding to the true model with $p$ parameters, as defined in (1), and we consider the set containing all possible auxiliary models and the true model, denoted as $A^* = A_{q_{max}} \cup \{(a^*, p)\}$. We then have the following proposition.

**Proposition 2** Given the models of Section 2 and assumptions A.1–A.10 in Appendix A, define the model $(\tilde{a}, \tilde{q})$ such that

$$(\tilde{a}, \tilde{q}) = \arg \min_{(q,a) \in A^*} IC_{LM}(a,q),$$

where

$$IC_{LM}(a,q) = -2 \sum_{i=1}^{N} \log \phi(x_{a,q,i}; \tilde{W}_{N,a,q}) + p(N,q), \quad (9)$$

where the penalty $p(N, q)$ satisfies assumption A.9. Then,

$$(\tilde{a}, \tilde{q}) \overset{P}{\to} (a^*, p), \quad as \ N,T \to \infty,$$

i.e. with probability one the criterion is asymptotically selecting the true model.\(^6\)

**Proof:** See Appendix B.

\(^6\)The previous probabilistic limit is a symbolic way of defining consistency. Indeed, $a$ is an index indicating a particular class of auxiliary models and not a true parameter. We could formalize it by considering a selecting vector of the same dimension as the number of models considered and taking values 1 or 0 if a given model is selected by the criterion or not, see Andrews (1999) and Hall et al. (2011). However, for simplicity of notation we prefer to avoid this choice as, once understood the meaning of $a$, nothing changes in the proofs.
The assumed form of the penalty is \( p(N, q) = h(q)K_N \), as postulated by Andrews (1999). In \( AIC_{IM}(a, q) \) we have \( h(q) = q(q+1)/2 \) and \( K_N = 1 \), hence \( p(N, q) = p(q) \) and it does not diverge with \( N \). Penalties similar to \( BIC \) and \( HQIC \) are potential candidates. Andrews (2000) and Hall et al. (2007) report simulation evidence that the \( BIC \)-type penalty –i.e. \( K_N = \log N \)– works best in their context, since \( K_N = \log \log N \) appears to be a too slow rate. This is also our experience. Bozdogan (1987) suggests \( K_N = (\log N)^b \) for \( b \geq 1 \) as a generalization of \( BIC \). The Monte Carlo study in the next section is performed with \( K_N = \log N \).

The criteria and other matching–based methods Other matching–based estimation methods, such as GMM, MSM and MSQ, also require the choice of functions. Provided that these functions ensure consistency of the estimates, all methods deliver asymptotically Gaussian estimates; their precision being the only difference. In GMM and MSM \( q \) is the number of moment conditions that provide consistent estimators. Then our criteria become similar to the RMSC of Hall et al. (2007). In MSQ the issue of weak identification is not relevant since, in theory, any \( q \)-dimensional vector of functions of quantiles provide consistent estimators. We can therefore, in principle, estimate the model of interest with different matching methods (nested and non–nested models, different moment conditions, and different functions of quantiles), and select the one that provides the smallest indirect criteria. A deeper understanding of the selection process between and within inference methods remains open and deserves further investigation.

4 Monte Carlo Study

To investigate the performance of the indirect criteria, we proceed with a thorough Monte Carlo study on two simple and widely used models: a moving average of order one and a stochastic volatility model. Appendix C shows the step–by–step description of the implementation.

Moving Average Consider the following MA(1) model

\[
y_t = \mu + \theta u_{t-1} + u_t, \quad t = 1 \ldots, T,
\]

where \( u_t \) is a standardized Gaussian white noise. The number of parameters is \( p = 2 \). We set \( \mu = 0.1 \) and we simulate \( R = 500 \) series for \( T = \{1000, 10000\} \) and \( \theta = \{0.2, 0.5, 0.8\} \). The purpose of this study is to investigate the performance of the criteria when choosing among the true and nested auxiliary models, for which we choose a set of nested autoregressive models with intercept:

\[
y_t = \phi_0 + \phi_1 y_{t-1} + \ldots + \phi_r y_{t-r} + \varepsilon_t,
\]

where \( \varepsilon_t \) is a standardized Gaussian white noise and \( r = \{1, 2, 3, 4, 5, 6, 7, 8, 9\} \). The sample size of the simulated series \( x_a,q,i \) is set to \( N = \{100, 500, 1000\} \), and \( H = \{1, 10\} \). The aim of varying \( N \) is to study the differences between the penalties. The number of parameters of the auxiliary models is given by \( q = 2 \) for the MA(1) and \( q = r + 1 \) for the AR(\( r \)).
Tables 1–3 and 4–6 in Appendix D report the criteria for $T = 1000$ and 10000 respectively. The column Rate shows the percentage of times (out of $R = 500$) that an auxiliary model is chosen, while the columns Mean and Var show the sample mean and variance of the $R = 500$ replications.

We extract five conclusions. First, regardless of $T$, $N$ and $H$, on average (column Mean) $IC_{IM}$ chooses the true model (numbers in bold) followed by an AR (numbers in bold and italics) with lags that increase with $\theta$. An increase in $\theta$ increases the discrepancy in the criteria between the MA(1) and the nested AR models. This may be due to the fact that a small persistence in the MA(1) process is easily captured by few AR lags, and that’s also why the criteria chooses the AR when $\theta$ is small. However, large values of $\theta$ require a large number of lags in the AR model. This effect becomes more apparent once $N$ increases. Second, and related with the previous conclusion, the criteria have a very small dispersion (column Var), and in the majority of cases the $IC_{IM}$ criterion chooses the true model (column Rate) for any value of $N$. This is not the case for $AIC_{IM}$. As $N$ increases the frequency at which $AIC_{IM}$ selects models others than the true one increases. Given these two findings, we conclude that the penalty with $K_N = \log N$ works better than with $K_N = 1$. Third, on average $IC_{IM}$ always chooses an auxiliary AR model with equal or smaller lags than the $AIC_{IM}$ does. This is in line with the simulation evidence reported in Andrews (1999) and Hall et al. (2007) for choosing consistent and non–redundant moment conditions respectively. Fourth, an increase in $H$ stabilizes the results and increases the quality of the criteria. This is particularly clear for $AIC_{IM}$ and is a natural effect of the increase in the precision of the estimates leading automatically to an increase in the values of the likelihoods, and to a decrease of the criteria. Also an increase in $T$ stabilizes the quality of the criteria, as expected from Proposition 2.

Stochastic Volatility Consider the following model:

$$y_t = e^{ht/2}u_t, \quad u_t \sim N(0,1),$$
$$h_t = \mu + \rho h_{t-1} + v_t, \quad v_t \sim N(0,\sigma^2), \quad t = 1\ldots,T.$$ 

We set $\mu = 0$, $\sigma^2 = 0.01$ and we simulate three series, each of size $T = 10000$, for three choices of $\rho = \{0.2,0.4,0.9\}$. Other choices of $T$ (smaller) are possible but the results for the MA(1) have shown that $T$ does not play a significant role. The first two values correspond to processes with mild volatility clustering, while the third is a common choice in the existing literature (e.g. Shephard (1996) and Monfardini (1998)) and close to values often obtained for daily financial returns. Following Broto and Ruiz (2004), we refrain from estimating $\mu$. Therefore, the number of parameters is $p = 2$. Similar to the MA(1), we choose $N = \{100, 500, 1000\}$ and $H = \{1, 10\}$. The main purpose of this study is to investigate the performance of the criteria when choosing among nested and non–nested auxiliary models. The first class of auxiliary models consists of ARCH($r$) models:

$$y_t = d_t^{1/2}z_t, \quad z_t \sim N(0,1),$$
$$d_t = \alpha_0 + \alpha_1 y_{t-1}^2 + \ldots + \alpha_r y_{t-r}^2,$$

for $r = \{1, \ldots, 9\}$. The number of auxiliary parameters is $q = r + 1$. This class of auxiliary models was proposed by Engle and Lee (1996) and Calzolari et al. (2008). We
also consider a GARCH(1,1) \( d_t = \alpha_0 + \alpha_1 y_{t-1}^2 + \alpha_2 d_{t-1} \) and hence \( q = 3 \). The second class consists of AR(\( r \)) models on the log of the squared observations:

\[
\ln(y_t^2) = \beta_0 + \beta_1 \ln(y_{t-1}^2) + \ldots + \beta_r \ln(y_{t-r}^2) + \eta_t, \quad \eta_t \sim N(0, \varphi)
\]

with \( r = \{1, \ldots, 9\} \). The number of auxiliary parameters is \( q = r + 2 \). This class of auxiliary models was proposed by Monfardini (1998). We also consider an ARMA(1,1)

\[
\ln(y_t^2) = \beta_0 + \beta_1 \ln(y_{t-1}^2) + \beta_2 \eta_{t-1} + \eta_t \quad \text{and hence} \quad q = 3.
\]

Tables 7–9 in Appendix D report the criteria. We extract three conclusions, all in line with the findings of the MA(1) model. First, similarly to the results reported by Monfardini (1998), the lag selected by the criteria, for both ARCH(\( r \)) and AR(\( r \)), increases with the persistence of the volatility. More importantly, regardless of the value of \( H, N \), and the penalty, the (G)ARCH model is always preferred (to a large extend) to the AR(MA), suggesting that the former is closer to the SV than the latter. Moreover, GARCH(1,1) is almost always preferred to an ARCH(\( r \)), specially for large values of \( \rho \). Second, increasing \( H \) stabilizes the results and decreases the values of the criteria. Third, between the ARCH(\( r \)) and AR(\( r \)) models, \( IC_{IM} \) always chooses a model equal or smaller than the \( AIC_{IM} \) does. This confirms once more that \( K_N = \log N \) works better than \( K_N = 1 \).

## 5 Conclusions

Indirect estimation methods depend on the choice of an auxiliary model. The difference between them comes in terms of precision of the estimators and, theoretically, the larger the number of parameters of the auxiliary model, the more precise are the estimates of the model of interest. But to account for the trade–off between the fit of the auxiliary model to the model of interest and its dimension, we need a penalty that depends on the number of parameters in the auxiliary model. In this article we develop information criteria for selecting among nested and non–nested auxiliary models. The criteria can also be used for choosing functions to match between and within other inference methods, such as GMM, the Method of Simulated Moments and the Method of Simulated Quantiles. We proceed with a thorough Monte Carlo study on a MA(1) and a stochastic volatility model. For the former we choose as auxiliary models a set of AR(\( r \)) and the true model. The criteria choose the true model as the best auxiliary model. As for the SV model, we consider two auxiliary classes of models: (G)ARCH and AR(MA) on the log of the square of the observations. Among the ARCH and AR type of models, the lag chosen by the criteria increases with the SV autoregressive parameter. Moreover, the criteria always prefer the (G)ARCH to the AR(MA) models.
References


A Assumptions

A.1 The set of parameters $\Theta$ is non-empty, compact, and $\theta_0 \in \text{int}(\Theta)$. The same holds for the auxiliary model: $B$ is non-empty, compact, and $\beta_0 \in \text{int}(B)$.

A.2 The log-likelihoods $\ell$ and $\ell_a$ are twice-differentiable, i.e. $\ell \in C^2(\Theta)$ and $\ell_a \in C^2(B)$.

A.3 The true parameter vector $\theta_0 \in \Theta$ is the unique maximum for $E[\ell(\theta; Y)]$ and $\beta_0$ is the unique maximum of $E[\ell_a(\beta; Y)]$. Moreover, $b(\theta_0)$ is the unique maximum of $E[\ell_a(b(\theta_0); Y)]$, which implies $\beta_0 = b(\theta_0)$.

A.4 The following matrices exist finite and are positive definite:

$$
I(\theta_0) = E_Y \left[ \frac{\partial \ell(\theta; Y)}{\partial \theta} \bigg|_{\theta = \theta_0} \left( \frac{\partial \ell(\theta; Y)}{\partial \theta'} \bigg|_{\theta = \theta_0} \right) \right],
$$

$$
I(\beta_0) = E_Y \left[ \frac{\partial \ell_a(\beta; Y)}{\partial \beta} \bigg|_{\beta = \beta_0} \left( \frac{\partial \ell_a(\beta; Y)}{\partial \beta'} \bigg|_{\beta = \beta_0} \right) \right] \quad \text{and}
$$

$$
J(\beta_0) = -E_Y \left[ \frac{\partial^2 \ell_a(\beta; Y)}{\partial \beta \partial \beta'} \bigg|_{\beta = \beta_0} \right].
$$

And the corresponding sample matrices converge in probability to them.

A.5 The binding function $b(\theta)$ is continuously differentiable, i.e. $b \in C^1(\Theta)$.

A.6 The binding function $b(\theta)$ is injective.

A.7 The $q \times p$ matrix $\frac{\partial b(\theta_0)}{\partial \theta}$ has full-column rank $p$.

A.8 $AIC_{1M}(a, q)$ has a unique minimum over $A_{\eta_{max}}$.

A.9 The penalty function is such that $p(N, q) = h(q)K_N$, with $h(q)$ strictly increasing and $K_N$ satisfying $K_N \to \infty$ and $K_N/N \to 0$ as $N \to \infty$.

A.10 The following matrices exist, are finite and positive definite:

$$
J_{a,q}(\hat{w}_{T,a,q}) = -E_{X_{a,q}} \left[ \frac{\partial^2 \log \phi(X_{a,q}; w_{a,q})}{\partial w_{a,q} \partial w_{a,q}'} \bigg|_{w_{a,q} = \hat{w}_{T,a,q}} \right],
$$

$$
J^*(\hat{w}_T^*) = -E_{X^*} \left[ \frac{\partial^2 \log \phi(X^*; w^*)}{\partial w^* \partial w^*'} \bigg|_{w^* = \hat{w}_T^*} \right].
$$

And the corresponding sample matrices converge in probability to them.

Assumptions A.1–A.4 are usual for ML estimation, and in a conditional case all log-likelihoods are to be intended as conditional on other variables and possibly on past information. Assumptions A.5–A.7 characterize the binding function and, in particular, A.6 and A.7 are the global and local identifiability conditions. A.8 and A.9 characterize the objective function of the information criterion. In particular they guarantee the existence of a global minimum and consistency respectively. Assumption A.10 is the analogous of A.4 but for the simulated data $X_{a,q}$ and $X^*$ used in the Propositions 1 and 2.
B  Technical appendix

We first define some quantities needed for the proofs, followed by two preliminary lemmas, and the proofs of Propositions 1 and 2.

Preliminary results

Definitions For simplicity of notation we omit, unless otherwise stated, all indexes referring to the class of auxiliary models \( a \), the size of the auxiliary parameter \( q \), and the sample size \( T \). We also omit any reference to the value of the parameters \( \theta_0 \) and \( \beta_0 \). We start with some definitions and then we prove two lemmas necessary to prove Proposition 1.

We re–write the asymptotic variance-covariance matrix given in (4) as

\[
W = (B' VB)^{-1},
\]

(10)

where \( V = J(\beta_0)I^{-1}(\beta_0)J(\beta_0) \) and \( B = \left. \frac{\partial h(\theta)}{\partial \theta} \right|_{\theta = \theta_0} \). Recall that \( V \) is \( q \times q \), \( B \) is \( q \times p \) with full column rank \( p \) (see assumption A7). We define the \( q \times p \) right–inverse matrix of \( B' \)

\[
(B')^+ = B(B'B)^{-1},
\]

(11)

such that \( B'(B')^+ = I_q \). Notice that this matrix always exists as long as \( B' \) has full column rank \( p \). Then, consider the \( q \times q \) matrix \( BB' \) and its eigenvalue decomposition \( BB' = P \Lambda P' \), where the columns of \( P \) are the eigenvectors \( (PP' = I_q) \) and \( \Lambda \) is a \( q \times q \) diagonal matrix of eigenvalues. Since from assumption A.6 we have \( q \geq p \), \( \Lambda \) has entries \( \lambda_i > 0 \) for \( i = 1, \ldots, p \) and \( \lambda_i = 0 \) for \( i = p + 1, \ldots, q \). Then, we define the generalized inverse matrix

\[
(BB')^+ = P \Lambda^+ P',
\]

(12)

where

\[
\Lambda^+ = \begin{pmatrix}
L^{-1} & 0_{p \times (q-p)} \\
0_{(q-p) \times p} & 0_{(q-p) \times (q-p)}
\end{pmatrix} \quad \text{and} \quad L^{-1} = \text{diag}(\lambda_1^{-1} \ldots \lambda_p^{-1}).
\]

(13)

Using these definitions we define a \( q \times p \) matrix

\[
B^1 = B'(BB')^+.
\]

(14)

Finally, for the \( q \times q \) matrix \( V^{-1} \), we define the duplication matrix, \( D_q \) of dimension \( q^2 \times \frac{q(q+1)}{2} \), while for the \( p \times p \) matrix \( W \), we define the duplication matrix \( D_p \) of dimension \( p^2 \times \frac{p(p+1)}{2} \) which has a left–inverse \( D_p^+ = (D_p'D_p)^{-1}D_p' \), such that \( D_p^+D_p = I_{p(p+1)} \). These matrices are such that

\[
\text{vec}(V^{-1}) = D_q \text{vech}(V^{-1}), \quad \text{vech}(W) = D_p^\top \text{vec}(W).
\]

(15)

The following Lemma decomposes the asymptotic variance–covariance matrix \( W \) into the binding function and the information matrix of the auxiliary model.

**Lemma 1** Let \( w = \text{vech}(W) \) be a \( \frac{p(p+1)}{2} \)–dimensional vector and \( W \) the asymptotic variance–covariance matrix (4). Then \( w = Av \), where \( A \) is a \( \frac{p(p+1)}{2} \times \frac{q(q+1)}{2} \) matrix defined as \( A = \)
$D_p^+(B^1 \otimes B^+)D_q$, and $v$ is a $\frac{q(q+1)}{2} \times 1$ vector defined as $v = \text{vech}(V^{-1})$.

**Proof.** From (10), we have

$$W = (B'VB)^{-1} = C_1V^{-1}C_2,$$

where $C_1$ and $C_2$ must be such that

$$B'VBC_1V^{-1}C_2 = I_p,$$  \hspace{1cm} (16)

and

$$C_1V^{-1}C_2B'VB = I_p.$$  \hspace{1cm} (17)

We first proof the equality in (16). By substituting in (16) $C_1 = B^1$ (see (14)) and $C_2 = (B')^+$ (see (11)), and using (12) and (13), we have

$$B'VB^1V^{-1}B^+ = B'VB'B(BB')^+V^{-1}B^+$$

$$= B'VPA'PA'^+P'V^{-1}B(B'B)^{-1}$$

$$= B'VP \left( I_p \begin{array}{c} 1_p \end{array} \right) \begin{array}{c} 0_p \end{array} \begin{array}{c} 1_{p \times (q-p)} \end{array} \begin{array}{c} 0_{(q-p) \times (q-p)} \end{array} \begin{array}{c} 1_{(q-p) \times (q-p)} \end{array} \right) P'V^{-1}B(B'B)^{-1}$$

$$= B'VP(p)p'TV^{-1}B(B'B)^{-1}$$

$$= B'V1V^{-1}B(B'B)^{-1}$$

$$= B'B(B'B)^{-1}$$

$$= I_p,$$

where $P^{(p)}$ is a $q \times p$ matrix with just the first $p$ columns of $P$. The equality in (17) is proven analogously.

Then

$$W = B'V^{-1}(B')^+ = B^1V^{-1}(B^+)^,'$$

and, from the properties of the vec$(\cdot)$ operator,

$$\text{vec}(W) = \text{vec}(B^1V^{-1}(B^+)^,) = (B^1 \otimes B^+) \text{vec}(V^{-1}).$$  \hspace{1cm} (18)

Finally, from (18), and using the duplication matrices defined in (15), we have

$$w = \text{vech}(W) = D_p^+\text{vec}(W)$$

$$= D_p^+ \left( B^1 \otimes B^+ \right) \text{vec}(V^{-1})$$

$$= D_p^+ \left( B^1 \otimes B^+ \right) D_q\text{vech}(V^{-1})$$

$$= A^pv.$$  

In order to estimate the binding function, we have to simulate $H$ paths $(y^1_1(\theta), \ldots, y^1_{J^1}(\theta))$ of length $T^*$ not necessarily equal to $T$. Thus, $b(\theta)$ is estimated using $\hat{\beta}_{HT^*}(\theta)$ as defined in (2). The error of this estimator can be made as small as we wish by taking $H \rightarrow \infty$ and $T^* \rightarrow \infty$. Moreover, we estimate derivatives numerically as

$$\frac{\partial \hat{\beta}_{HT^*}(\theta)}{\partial \theta} \sim \frac{-\hat{\beta}_{HT^*}(\theta + 2\Delta) + 8\hat{\beta}_{HT^*}(\theta + \Delta) - 8\hat{\beta}_{HT^*}(\theta - \Delta) + \hat{\beta}_{HT^*}(\theta - 2\Delta)}{12\Delta}.$$
for some arbitrarily small $\Delta$. Thus,

$$
\lim_{H,T^* \to \infty} \left( \lim_{\Delta \to 0} \frac{\partial \hat{b}_{H,T^*}(\theta)}{\partial \theta} \right) = \frac{\partial b(\theta)}{\partial \theta} = B.
$$

Since $\Delta$, $H$, and $T^*$ can be arbitrarily fixed, in simulations we can neglect the error made in estimating $B$ and thus consider $A$, which involves $B$, as non-random. Intuitively, $A$ does not involve any additional information with respect to the information contained in the elements of the asymptotic variance–covariance matrix of the auxiliary model $V^{-1}$.

We now re-establish the complete notation, and the estimator of the asymptotic variance–covariance matrix is such that

$$
\hat{w}_{T,a,q} = A \hat{v}_{T,a,q}.
$$

Recall that for the simulated vector $X_{a,q} \equiv \sqrt{T}(\hat{\theta}_{T,IM,a,q} - \theta_0)$ we have a sample covariance matrix $\tilde{w}_{N,a,q}$. If we define a new vector $Y_{a,q} \equiv \sqrt{T}(\beta_{T,IM,a,q} - \beta_0)$, then the sample covariance matrix of $Y_{a,q}$ is $\tilde{v}_{N,a,q}$ such that

$$
\tilde{w}_{N,a,q} = A \tilde{v}_{N,a,q}.
$$

The following Lemma shows the asymptotic distribution of $\tilde{v}_{N,a,q}$.

Lemma 2 For any given auxiliary model $a$ with parameters of dimension $q$, we have, for $N \to \infty$,

$$
\sqrt{N}(\tilde{v}_{N,a,q} - \hat{v}_{T,a,q}) \sim N\left(0, J_{a,q}^{-1}(\hat{w}_{T,a,q})A^\dagger\right),
$$

where $N$ is the size of the simulated process $(x_{a,q},1,\ldots,x_{a,q},N)$,

$$
J_{a,q}(\hat{w}_{T,a,q}) = -E_{x_{a,q}} \left[ \frac{\partial^2 \log \phi(X_{a,q}, w_{a,q})}{\partial w_{a,q} \partial w'_{a,q}} \right]_{w_{a,q} = \hat{w}_{T,a,q}}.
$$

and $A^\dagger = (A'A)^+A$ with $(A'A)^+$ defined analogously as in (12), but this time with $\frac{p(p+1)}{2}$ non-zero eigenvalues, thus $A^\dagger A = I_{q(q+1)}$.

Proof. We defined $\tilde{w}_{N,a,q}$ to be the vector containing the entries of the sample covariance matrix of $X_{a,q}$ which in turn is distributed as a Gaussian with zero mean and covariance matrix $\overline{W}_{T,a,q}$. Then by defining $\hat{w}_{T,a,q} = \text{vech}(\overline{W}_{T,a,q})$, we have, as $N \to \infty$,

$$
\sqrt{N}(\tilde{w}_{N,a,q} - \hat{w}_{T,a,q}) \sim N\left(0, J_{a,q}^{-1}(\hat{w}_{T,a,q})\right).
$$

Using Lemma 1, the estimator in (19), and the delta method, we obtain the result. □

Proof of Proposition 1

In this and the following proofs, we reintroduce those indexes referring to the class of auxiliary models $a$, the size of the auxiliary parameter $q$, and the observed and simulated samples sizes $T$ and $N$. We also indicate the value of the parameters in which the matrices are computed, but we simplify the notation using the following definitions of the indirect methods and ML asymptotic
If we replace the expectation by its sample counterpart, we have to minimize

\[ \theta \]

Finally, notice that, since ML estimation of the model of interest is unfeasible, we cannot compute the form, generically denoted as \( W \).

Moreover, analogously to the \( W \) we define the \( W \), respectively. These matrices are the consistent ML estimators of \( \hat{W}_{T,a,q} \) and \( \hat{W}_T \) respectively. For any of the matrices \( W \) defined above, we have a vectorized form, generically denoted as \( w \equiv \text{vech}(W) \).

Finally, notice that, since ML estimation of the model of interest is unfeasible, we cannot compute \( \theta_{T,ML} \) and therefore \( X^* \) cannot be simulated. This, however, is an object needed only for the proof and plays no role in the final formulation of the criteria.

Given consistency of the sample variance–covariance matrices we have, as \( N \to \infty \),

\[ \sqrt{N} (\hat{w}_{N,a,q} - \bar{w}_{T,a,q}) \sim N(0, J_{a,q}^{-1}(\bar{w}_{T,a,q})) \]  

and

\[ \sqrt{N} (\hat{w}^*_N - \bar{w}^*_T) \sim N(0, J^*^{-1}(\bar{w}^*_T)) \],

where the asymptotic variance–covariance matrices have now size \( \frac{p(p+1)}{2} \times \frac{p(p+1)}{2} \) and are defined as

\[ J_{a,q}(\bar{w}_{T,a,q}) = -E_{X_{a,q}} \left[ \frac{\partial^2 \log \phi(X_{a,q}; w_{a,q})}{\partial w_{a,q} \partial w_{a,q}'} \bigg|_{w_{a,q} = \bar{w}_{T,a,q}} \right] \] and

\[ J^*(\bar{w}^*_T) = -E_{X^*} \left[ \frac{\partial^2 \log \phi(X^*; w^*)}{\partial w^* \partial w^*'} \bigg|_{w^* = \bar{w}^*_T} \right]. \]

The sample analogues of the two previous expressions are:

\[ \tilde{J}_{N,a,q}(\bar{w}_{T,a,q}) = -\frac{1}{N} \sum_{i=1}^{N} \frac{\partial^2 \log \phi(X_{a,q,i}; w_{a,q})}{\partial w_{a,q} \partial w_{a,q}'} \bigg|_{w_{a,q} = \bar{w}_{T,a,q}} \] and

\[ \tilde{J}^*(\bar{w}^*_T) = -\frac{1}{N} \sum_{i=1}^{N} \frac{\partial^2 \log \phi(X_{a,q,i}^*; w^*)}{\partial w^* \partial w^*'} \bigg|_{w^* = \bar{w}^*_T}. \]

Once \( X_{a,q} \) is given, we have to find an auxiliary model with asymptotic variance–covariance matrix \( W_{a,q} \) that better approximates (in terms of information) the best achievable value \( W_0 \). In order to do this, we minimize the Kullback–Leibler information (or negentropy)

\[ D(w_0^*, w_{a,q}) = 2 E_{X^*} \left[ \log \frac{\phi(X^*; w_0^*)}{\phi(X_{a,q}; w_{a,q})} \right], \]

If we replace the expectation by its sample counterpart, we have to minimize

\[ \tilde{D}(w_0^*, w_{a,q}) = \frac{2}{N} \sum_{i=1}^{N} \log \frac{\phi(X_{a,q,i}^*; w_0^*)}{\phi(X_{a,q,i}; w_{a,q})}. \]
Since we can only estimate \( w_{a,q} \), the best we can do is to use its indirect estimator \( \tilde{w}_{T,a,q} \). Therefore, instead of minimizing (24) we look for the auxiliary model defined by the couple \((a, q)\) that minimizes

\[
E_w \left[ \hat{D}_N(w^*_0, \tilde{w}_{T,a,q}) \right] = E_w \left[ \frac{2}{N} \sum_{i=1}^{N} \log \frac{\phi(x_i^*; w^*_0)}{\phi(x_{a,q,i}; \tilde{w}_{T,a,q})} \right].
\]

(25)

We decompose the objective function as the sum of two terms

\[
Q \equiv \hat{D}_N(w^*_0, \tilde{w}_{T,a,q}) = \frac{\hat{D}_N(w^*_0, \tilde{w}_{N,a,q}) + \hat{D}_N(\tilde{w}_{N,a,q}, \tilde{w}_{T,a,q})}{g}.
\]

(26)

Term \( P \) of (26) can be written as

\[
P = \hat{D}_N(w^*_0, \tilde{w}_{N,a,q}) = \hat{D}_N(w^*_0, \tilde{w}^*_{N,a,q}) + \hat{D}_N(\tilde{w}^*_N, \tilde{w}_{N,a,q}) + \hat{D}_N(\tilde{w}_{N,a,q}, \tilde{w}_{T,a,q}).
\]

So (26) can be written as

\[
Q = A + C + B + G.
\]

(27)

Both term \( A \) and \( C \) do not depend on the chosen auxiliary model thus it can be treated as constants in the criterion. Let us consider the other two terms \( B \) and \( G \) separately.

\( B \): This term is

\[
B = \frac{2}{N} \sum_{i=1}^{N} \log \frac{\phi(x_i^*; \tilde{w}^*_N)}{\phi(x_{a,q,i}; \tilde{w}_{N,a,q})} = \frac{2}{N} \sum_{i=1}^{N} \left\{ \log \phi(x_i^*; \tilde{w}^*_N) - \log \phi(x_{a,q,i}; \tilde{w}_{N,a,q}) \right\}.
\]

Only the second term on the right hand side depends on the chosen auxiliary model, \((a, q)\), while the first term can be treated as a constant.

\( G \): This term can be approximated using a Taylor expansion in a neighborhood of \( \tilde{w}_{N,a,q} \):

\[
G \approx -(\tilde{w}_{T,a,q} - \tilde{w}_{N,a,q})' \frac{1}{N} \sum_{i=1}^{N} \frac{\partial^2 \log \phi(x_{i,a,q}; w_{a,q})}{\partial w_{a,q} \partial w_{a,q}'} \bigg|_{w_{a,q}=\tilde{w}_{N,a,q}} \tilde{w}_{T,a,q} - \tilde{w}_{N,a,q} = \tilde{w}_{T,a,q} - \tilde{w}_{N,a,q}.
\]

\[
J_{N,a,q}(\tilde{w}_{N,a,q}) = J_{a,q}(\tilde{w}_{T,a,q}).
\]

Since

\[
\lim_{N \to \infty} \frac{J_{N,a,q}(\tilde{w}_{N,a,q}) - J_{a,q}(\tilde{w}_{T,a,q})}{N} = 0,
\]

and given (21) and lemma 2, we have for \( N \to \infty \)

\[
NG \approx N(\tilde{w}_{T,a,q} - \tilde{w}_{N,a,q})' J_{a,q}(\tilde{w}_{T,a,q}) (\tilde{w}_{T,a,q} - \tilde{w}_{N,a,q}) = N(\tilde{w}_{T,a,q} - \tilde{w}_{N,a,q})' A^T J_{a,q}(\tilde{w}_{T,a,q}) A (\tilde{w}_{T,a,q} - \tilde{w}_{N,a,q}) \sim \chi^2_{q(q+1)/2}.
\]

By combining the previous results, we have that minimizing (25) is equivalent to minimize

\[
E_w [NQ] = E_w [NA + NC + NB + NG] = k - 2 \sum_{i=1}^{N} \log \phi(x_{a,q,i}; \tilde{w}_{N,a,q}) + \frac{q(q+1)}{2}.
\]
where the first term $k$ on the right hand side contains $A$, $C$, and the first term of $B$, therefore it does not depend on the chosen auxiliary model, $(a, q)$. Thus, minimizing (25) is equivalent to minimize with respect to all possible auxiliary models $(a, q)$ the following criterion

$$AIC_{IM}(a, q) = -2 \sum_{i=1}^{N} \log \phi(x_{a,q,i}; \tilde{w}_{N,a,q}) + \frac{q(q+1)}{2},$$

where $\tilde{w}_{N,a,q}$ is the vector containing the entries of the sample variance–covariance matrix of the simulated vector $X_{a,q}$.

**Proof of Proposition 2**

For each model $(a, q) \in \mathcal{A}^*$, we have the criterion

$$IC_{IM}(a, q) = -2 \sum_{i=1}^{N} \log \phi(x_{a,q,i}; \tilde{w}_{N,a,q}) + h(q)K_N,$$

If we could compute the criterion in correspondence of the true model, we would have

$$IC_{IM}(a^*, p) = -2 \sum_{i=1}^{N} \log \phi(x_{a,q,i}^*; \tilde{w}_{N,a,q}^*) + h(p)K_N.$$ 

Therefore,

$$\frac{1}{N}[IC_{IM}(a, q) - IC_{IM}(a^*, p)] = \frac{2}{N} \sum_{i=1}^{N} \frac{\phi(x_{a,q,i}^*; \tilde{w}_{N,a,q})}{\phi(x_{a,q,i}; \tilde{w}_{N,a,q})} + (h(q) - h(p))\frac{K_N}{N} = \tilde{D}_N(\tilde{w}_N^*, \tilde{w}_{N,a,q}) + (h(q) - h(p))\frac{K_N}{N}. \quad (28)$$

The first term on the right–hand–side of (28) can be decomposed as

$$\tilde{D}_N(\tilde{w}_N^*, \tilde{w}_{N,a,q}) = \tilde{D}_N(\tilde{w}_N^*, \tilde{w}_{T,a,q}^*) + \tilde{D}_N(\tilde{w}_{T,a,q}^*, \tilde{w}_0^*) + \tilde{D}_N(\tilde{w}_0^*, \tilde{w}_{0,a,q}) + \tilde{D}_N(\tilde{w}_{0,a,q}, \tilde{w}_{T,a,q}) + \tilde{D}_N(\tilde{w}_{T,a,q}, \tilde{w}_{N,a,q}). \quad (29)$$

From (21) and (22), and using Taylor expansions, we have that as $N \to \infty$ the first and last term of (29) are $\mathcal{O}_p(N^{-1})$. In a similar way, using Taylor approximations we write

$$\tilde{D}_N(\tilde{w}_0^*, \tilde{w}_{T,a,q}) \simeq (\tilde{w}_0^* - \tilde{w}_{T,a,q}^*)/\tilde{J}_{T,a,q}(\tilde{w}_{T,a,q}^*)(\tilde{w}_0^* - \tilde{w}_{T,a,q}^*).$$

From Gouriéroux et al. (1993) we have, as $T \to \infty$,

$$p \lim_{T \to \infty} \tilde{w}_{T,a,q} = \tilde{w}_{0,a,q},$$

which, provided that $\tilde{J}_{T,a,q}$ is bounded by A.10, implies $\|\tilde{w}_0^* - \tilde{w}_{T,a,q}^*\| = \mathcal{O}_p(T^{-1/2})$ and therefore the fourth term of (29) is $\mathcal{O}_p(T^{-1})$. Analogously from Maximum Likelihood theory we have that the second term in (29) can be approximated as

$$\tilde{D}_N(\tilde{w}_0^* - \tilde{w}_T^*),$$
which, given consistency of \( \hat{\omega}_T^* \) as \( T \to \infty \), and boundedness of \( \hat{J}_T^* \) (see A.10), is also \( O_p(T^{-1}) \).

Therefore, (28) becomes

\[
\frac{1}{N}[IC_{IM}(a, q) - IC_{IM}(a^*, p)] = \hat{D}_N(w_0^*, w_{0,a,q}) + O_p(N^{-1}) + O_p(T^{-1}) + (h(q) - h(p)) \frac{K_N}{N}.
\]

And, since \( K_N/N = o(1) \),

\[
p \lim_{N \to \infty} \left\{ \frac{1}{N}[IC_{IM}(a, q) - IC_{IM}(a^*, p)] \right\} = p \lim_{N \to \infty} \left\{ \hat{D}_N(w_0^*, w_{0,a,q}) \right\} + O_p(T^{-1}) = D(w_0^*, w_{0,a,q}) + O_p(T^{-1}).
\]

If we also let \( T \to \infty \), we have

\[
p \lim_{N,T \to \infty} \left\{ \frac{1}{N}[IC_{IM}(a, q) - IC_{IM}(a^*, p)] \right\} = D(w_0^*, w_{0,a,q}) \geq 0,
\]

where the last inequality follows from the definition of Kullback–Leibler information which would be zero when \( w_{0,a,q} = w_0^* \).

From (30), we have that, as \( N, T \to \infty \),

\[
IC_{IM}(a^*, p) \leq IC_{IM}(a, q), \quad \text{for any } (a, q) \in \mathcal{A}^*,
\]

with probability one. Moreover, given the definition of \( (\tilde{a}, \tilde{q}) \) and A.8, we also have

\[
IC_{IM}(\tilde{a}, \tilde{q}) < IC_{IM}(a, q), \quad \text{for any } (a, q) \in \mathcal{A}_{q_{\text{max}}}.
\]

Therefore, as \( N, T \to \infty \), we have that with probability one and when considering all models in \( \mathcal{A}^* \), the models \( (\tilde{a}, \tilde{q}) \) and \( (a^*, p) \) will coincide. In other words we would select the true model if it is included in the set of models we are considering. □
C  Step–by–step implementation

Step 1: Choose an auxiliary model $M^a(y, \beta)$ and estimate $\beta$ from the observations to obtain $\hat{\beta}_T$.

Step 2: Find the Indirect Inference estimator of $\theta$:

$$\hat{\theta}_{HT,IM,a,q} = \arg\min_{\theta \in \Theta} \left( \hat{\beta}_{HT}(\theta) - \hat{\beta}_T \right)' \Omega^* \left( \hat{\beta}_{HT}(\theta) - \hat{\beta}_T \right),$$

where $\Omega^*$ is the optimal weighting matrix. The algorithm for the minimization starts at some initial value $\theta^0$. The corresponding simulated $H$ paths of length $T$ are $y^1_H(\theta^0), \ldots, y^H_H(\theta^0)$, $h = 1, \ldots, H$, and the estimates $\hat{\beta}_{HT}(\theta^0)$. Note that the simulated paths can be of size $T^* > T$.

Step 3: Let $X_{a,q} \equiv \sqrt{T}(\hat{\theta}_{HT,IM,a,q} - \theta_0)$. Simulate a sample of length $N$ from the $N(0, \hat{\theta}_{HT,IM,a,q}(\hat{\theta}_{HT,IM,a,q}))$. Denote the $i$–th simulated observation of $X_{a,q}$ as the $p$–dimensional vector $x_{a,q,i} = (x_{1,a,q,i} \ldots x_{p,a,q,i})'$.

Step 5: From the $N$ simulated series $x_{a,q,i}$ compute the sample variance–covariance matrix $\hat{\theta}_{N,a,q}$ and the preferred information criterion given in (5) or (9).

Step 6: Select another (nested or non–nested) auxiliary model and repeat steps 1–5.

Step 7: Choose the auxiliary model that provides the smallest value for the preferred information criterion.
The model of interest is a $MA(1)$ with intercept $0.1$, moving average parameter $0.2$ and variance of the errors equal to $1$. The total number of replications is $R = 500$ and the simulated time series contains $T = 1000$ observations. The number of simulated paths $H$ is $1, 10, 500$ and $1000$. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.

### Table 1: $MA(1)$ with $\theta = 0.2$ and $T = 1000$

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<th>$H = 500$</th>
<th>$H = 1000$</th>
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<td>Rate Mean Var</td>
<td>Rate Mean Var</td>
<td>Rate Mean Var</td>
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<td>0.0015</td>
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The model of interest is a $MA(1)$ with intercept $0.1$, moving average parameter $0.2$ and variance of the errors equal to $1$. The total number of replications is $R = 500$ and the simulated time series contains $T = 1000$ observations. The number of simulated paths $H$ is $1$ and $10$, and $N$ (first column) equals $100, 500, 1000$. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.
The model of interest is a $MA(1)$ with intercept $\theta = 0.1$, moving average parameter $0.5$ and variance of the errors equal to $1$. The total number of replications is $R = 500$ and the simulated time series contains $T = 1000$ observations. The number of simulated paths $H$ is $1$ and $10$, and $N$ (first column) equals $100$, $500$, and $1000$. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.

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<td>$IC_{1M}$</td>
</tr>
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<td>Mean</td>
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Table 2: $MA(1)$ with $\theta = 0.5$ and $T = 1000$
Table 3: $MA(1)$ with $\theta = 0.8$ and $T = 1000$

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<th>$IC_{1M}$ Rate Mean Var</th>
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<tr>
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<tr>
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<td>$AR(8)$</td>
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<td>0 21.0540 0.0008</td>
<td>0 18.1354 0.0007</td>
<td>0 19.9813 0.0008</td>
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The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.8 and variance of the errors equal to 1. The total number of replications is $R = 500$ and the simulated time series contains $T = 1000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.
The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.2 and variance of the errors equal to 1. The total number of replications is $R = 500$ and the simulated time series contains $T = 10000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.

Table 4: $MA(1)$ with $\theta = 0.2$ and $T = 10000$

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The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.2 and variance of the errors equal to 1. The total number of replications is $R = 500$ and the simulated time series contains $T = 10000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.
The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.5 and variance of the errors equal to 1. The total number of replications is $R = 500$ and the simulated time series contains $T = 10000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.

Table 5: $MA(1)$ with $\theta = 0.5$ and $T = 10000$

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The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.5 and variance of the errors equal to 1. The total number of replications is $R = 500$ and the simulated time series contains $T = 10000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.
Table 6: $MA(1)$ with $\theta = 0.8$ and $T = 10000$

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The model of interest is a $MA(1)$ with intercept 0.1, moving average parameter 0.8 and variance of the errors equal to 1. The total number of replications is $H = 500$ and the simulated time series contains $T = 10000$ observations. The number of simulated paths $H$ is 1 and 10, and $N$ (first column) equals 100, 500, and 1000. The auxiliary model is either the model of interest itself or an $AR(r)$ (column Aux. Model). The table reports the percentage of choosing each model as the best (column Rate), the average value and the variance of the criteria (columns Mean and Var respectively) over the 500 replications. Numbers in bold denote the model chosen by the criteria. Numbers in bold and italics denote the second best model chosen by the criteria.
The model of interest is a SV with intercept 0, autoregressive parameter 0.2, and variances of the errors 1 and 0.01 for the observation and volatility equations respectively. The number of simulated paths \( H \) is 1 and 10, \( N \) (first column) equals 100, 500, and 1000, and \( T = 10000 \). The auxiliary models (column Aux. Model) are ARCH\( (r) \) and GARCH\( (1,1) \) on the observations (top three panels), and AR\( (r) \) and ARMA\( (1,1) \) in the log of the square observations (bottom three panels). Numbers in bold denote the model chosen by the criteria.

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The model of interest is a SV with intercept 0, autoregressive parameter 0.4, and variances of the errors 1 and $0.01$ for the observation and volatility equations respectively. The number of simulated paths $H$ is 1 and 10, $N$ (first column) equals 100, 500, and 1000, and $T = 10000$. The auxiliary models (column Aux. Model) are ARCH($r$) and GARCH(1,1) on the observations (top three panels), and AR($r$) and ARMA(1,1) in the log of the square observations (bottom three panels). Numbers in bold denote the model chosen by the criteria.

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Table 8: SV with $\rho = 0.4$.
Table 9: SV with $\rho = 0.9$

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The model of interest is a SV with intercept 0, autoregressive parameter 0.9, and variances of the errors 1 and 0.01 for the observation and volatility equations respectively. The number of simulated paths $H$ is 1 and 10, $N$ (first column) equals 100, 500, and 1000, and $T = 10000$. The auxiliary models (column Aux. Model) are $ARCH(r)$ and $GARCH(1,1)$ on the observations (top three panels), and $AR(r)$ and $ARMA(1,1)$ in the log of the square observations (bottom three panels). Numbers in bold denote the model chosen by the criteria.
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