Evaluation of Dynamic Stochastic General Equilibrium Models Based on Distributional Comparison of Simulated and Historical Data

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Abstract

We take as a starting point the existence of a joint distribution implied by different dynamic stochastic general equilibrium (DSGE) models, all of which are potentially misspecified. Our objective is to compare “true” joint distributions with ones generated by given DSGEs. This is accomplished via comparison of the empirical joint distributions (or confidence intervals) of historical and simulated time series. The tool draws on recent advances in the theory of the bootstrap, Kolmogorov type testing, and other work on the evaluation of DSGEs, aimed at comparing the second order properties of historical and simulated time series. We begin by fixing a given model as the “benchmark” model, against which all “alternative” models are to be compared. We then test whether at least one of the alternative models provides a more “accurate” approximation to the true cumulative distribution than does the benchmark model, where accuracy is measured in terms of distributional square error. Bootstrap critical values are discussed, and an illustrative example is given, in which it is shown that alternative versions of a standard DSGE model in which calibrated parameters are allowed to vary slightly perform equally well. On the other hand, there are stark differences between models when the shocks driving the models are assigned non-plausible variances and/or distributional assumptions.

\textit{JEL classification:} C12, C22.

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1 Introduction

In this paper, we merge recent econometric advances in bootstrapping time series and Kolmogorov type testing with recent developments in the evaluation of dynamic stochastic general equilibrium (DSGE) models. This is accomplished via the construction of a new tool for comparing the empirical joint distribution of historical time series with the empirical distribution of simulated time series.

Since the seminal papers by Kydland and Prescott (1982), Long and Plosser (1983) and King, Plosser and Rebelo (KPR: 1988a, b), there has been substantial attention given to the problem of reconciling the dynamic properties of data simulated from DSGE, and in particular from real business cycle (RBC) models, with the historical record. A partial list of advances in this area includes: (i) the examination of how RBC simulated data reproduce the covariance and autocorrelation functions of actual time series (see e.g. Watson (1993)); (ii) the comparison of DSGE and historical spectral densities (see e.g. Diebold, Ohanian and Berkowitz (1998)); (iii) the evaluation of the difference between the second order time series properties of vector autoregression (VAR) predictions and out-of-sample predictions from DSGE models (see e.g. Schmitt-Grohe (2000)); (iv) the construction of Bayesian odds ratios for comparing DSGE models with unrestricted VAR models (see e.g. Chang, Gomes and Schorfheide (2002), and Fernandez-Villaverde and Rubio-Ramirez (2004)); (v) the comparison of historical and simulated data impulse response functions (see e.g. Cogley and Nason (1995)); (vi) the formulation of “reality” bounds for measuring how close the density of an DSGE model is to the density associated with an unrestricted VAR model (see e.g. Bierens and Swanson (2000)); and (vii) loss function based evaluation of DSGE models (Schorfheide (2000)). The papers listed above are mainly concerned with the issue of model evaluation. Another strand of the literature is instead mainly concerned with providing alternatives to calibration (see e.g. De Jong, Ingram and Whiteman (2000) for a Bayesian perspective in which prior distributions are constructed around calibrated structural parameters). In most of the above papers, the issue of singularity (i.e. when the number of variables in the model is larger than the number of shocks) is circumvented by considering only a subset of variables, for which a nondegenerate distribution exists.\footnote{A novel alternative to calibration is proposed by Bierens (2005). He solves the singularity problem by considering a subset of variables for which a nondegenerate distribution exists.} Our work is closest to the first strand
of literature. In particular, our paper attempts to add to the model evaluation literature by introducing a measure of the "goodness of fit" of RBC models that is based on applying standard notions of Kolmogorov distance and drawing on advances in the theory of the bootstrap.²

The papers cited above primarily address the case in which the objective is to test for the correct specification of some aspects of a given candidate model. In the case of DSGE models, however, it is usually crucial to account for the fact that all models are approximations, and so are misspecified. This is the reason why the testing procedure that we develop allows us to evaluate the relative degree of misspecification of a given group of competing models based on the comparison of empirical distribution functions of historical data with those of DSGE simulated data. The DSGE models of interest in our context are simulated using both calibrated parameters (with calibrated values suggested by KPR, for example), and parameters estimated by using actual data, along the lines of Christiano (1988) and Christiano and Eichenbaum (1992). One important feature of our approach is that we begin by fixing a given DSGE model as the "benchmark" model, against which all "alternative" models are compared. The comparison is done using a distributional generalization of White's (2000) reality check, which assesses whether at least one of the alternative models provides a more "accurate" approximation to the true cumulative distribution than does the benchmark model. One key element of our approach is that we measure "accuracy" in terms of square error, as in Corradi and Swanson (2005a). We also outline the relationship between our measure of accuracy and the Kullback-Leibler Information Criterion (KLIC). DSGE model evaluation based on KLIC measures of accuracy is considered by Fernandez-Villaverde and Rubio-Ramirez (2004), and Chan, Gomes and Schorfheide (2002). For extensions of the methodology in this paper to the case of predictive density and conditional confidence interval accuracy evaluation, the reader is referred to Corradi and Swanson (2005b,c).

²In recent years, Kolmogorov type distance measures for testing distributional equality have been extended to the case of dependent observations and/or parameter estimation error (see e.g. Andrews (1997), Bai (2003)). Tests of this sort generally have limiting distributions that are not nuisance parameter free, and critical values cannot be tabulated. Papers addressing this issue in the bootstrap literature include Andrews (2002), Goncalves and White (2004), Hall and Horowitz (1996), Horowitz (2003), Inoue and Shintani (2005), Naik-Nimbalkar and Rajarshi (1994).
As mentioned above, our statistic is based on comparison of historical and simulated distributions. The limiting distribution of the statistic is a functional over a Gaussian process with a covariance kernel that reflects the contribution of parameter estimation error. This limiting distribution is thus not nuisance parameter free, and critical values cannot be tabulated. In order to obtain valid asymptotic critical values, we suggest two block bootstrap procedures, each of which depends on the relative rate of growth of the actual and simulated sample size. In addition, we circumvent the issue of singularity by considering a subset of variables (and their lagged values) for which a non singular distribution exists.

Our testing framework can be used to address questions of the following sort: (i) For a given DSGE model, what is the relative usefulness of different sets of calibrated parameters for mimicking different dynamic features of output growth? (ii) Given a fixed set of calibrated parameters, what is the relative performance of DSGE models driven by shocks with a different marginal distribution?

In order to illustrate how the proposed testing framework can be used, we consider the RBC model of Christiano (1988), characterized by flexible labor supply, capital depreciation, and two shocks - a permanent shock affecting technology and a transitory shock affecting preferences. Data are then simulated and various versions of the model are compared, in terms of their ability to reproduce the joint distribution of current output, lagged output, current hours worked, and lagged hours worked. The illustrations suggest that the methodology outlined in this paper provides a useful additional tool for examining the relevance of different RBC models vis a vis how well competing models capture the dynamic structural characteristics of the historical record.

The rest of the paper is organized as follows. Section 2 outlines the testing framework, describes the test statistic, and shows that the limiting distribution of the statistic is a zero mean Gaussian process with a covariance kernel that reflects both the contribution of parameter estimation error and the time series structure of the data. This is all done under the assumption that all models may be misspecified. In Section 3, the construction of bootstrap critical values is outlined, and the first order validity of the block bootstrap is established under two different assumptions with respect to the limit of the ratio of the actual sample size and the simulated sample period. An empirical illustration is given in Section 4, and concluding remarks are gathered in Section 5. All proofs are collected in the
2 Testing Framework

Our objective is to compare the joint distribution of historical variables with the joint distribution of simulated variables. Hereafter, for sake of simplicity, but without loss of generality, we limit our attention to the joint empirical distribution of (actual and model-based) current and previous period output. In principle, if we have a model driven by \( k \) shocks, then we can consider the joint CDF of \( k \) variables plus an arbitrary (but finite) number of lags. Consider \( m \) RBC models, and set model 1 as the benchmark model. Let \( \Delta \log X_t, t = 1, ..., T \) denote actual historical output (growth rates) and let \( \Delta \log X_{j,n}, j = 1, ..., m \) and \( n = 1, ..., S \), denote the output series simulated under model \( j \), where \( S \) denotes the simulated sample length. As mentioned above, some parameters are assumed to be kept fixed (at their calibrated values), while other parameters are estimated. Along these lines, denote \( \Delta \log X_{j,n}(\hat{\theta}_{j,T}) \), \( n = 1, ..., S \), \( j = 1, ..., m \) to be a sample of length \( S \) drawn (simulated) from model \( j \) and evaluated at the parameters estimated under model \( j \), where parameter estimation is done using the \( T \) available historical observations.\(^3\)

Now, let \( Y_t = (\Delta \log X_t, \Delta \log X_{t-1}) \) and \( Y_{j,n}(\hat{\theta}_{j,T}) = (\Delta \log X_{j,n}(\hat{\theta}_{j,T}), \Delta \log X_{j,n-1}(\hat{\theta}_{j,T})) \). Also, let \( F_0(u; \theta_0) \) denote the distribution of \( Y_t \) evaluated at \( u \) and \( F_j(u; \theta_j) \) denote the distribution of \( Y_{j,n}(\theta_j) \), where \( \theta_j \) is the probability limit of \( \hat{\theta}_{j,T} \), taken as \( T \to \infty \), and where \( u \in U \subset \mathbb{R}^2 \), possibly unbounded. Accuracy is measured in terms of the squared (approximation) error associated with model \( j, j = 1, ..., m \), defined as a (weighted) average over \( U \) of \( E \left( \left( F_j(u; \theta_j) - F_0(u; \theta_0) \right)^2 \right) \). Thus, the rule is to choose Model 1 over Model 2, say, if

\[
\int_U E \left( \left( F_1(u; \theta_j) - F_0(u; \theta_0) \right)^2 \right) \phi(u) du < \int_U E \left( \left( F_2(u; \theta_j) - F_0(u; \theta_0) \right)^2 \right) \phi(u) du,
\]

where \( \int_U \phi(u) du = 1 \) and \( \phi(u) \geq 0 \) for all \( u \in U \subset \mathbb{R}^2 \). For any evaluation point, this measure defines a norm and is a typical goodness of fit measure.

Another measure of distributional accuracy available in the literature (see e.g. Vuong (1989)), is the KLIC, according to which we should choose Model 1 over Model 2 if \( E(\log f_1(Y_t; \theta_1)) - \)

\(^3\)The reason why we use differences in the above is that the RBC model in our empirical illustration is characterized by permanent shocks to technology. In general, we require both the actual and the simulated series to be (strictly) stationary and mixing, and hence differencing may or may not be appropriate.
The KLIC thus chooses the model which on average gives higher probability to events which have actually occurred. Also, it leads to simple Likelihood Ratio tests. Our approach is an alternative to the KLIC that should be viewed as complementary in some cases, and preferred in others. For example, if we are interested in measuring accuracy for a given confidence interval, this cannot be done in an obvious manner using the KLIC, while it can easily be done using our measure. Furthermore, we often do not have an explicit form for the density implied by the various models we are comparing. Of course, model comparison could then be done using kernel density estimators, at the cost of nonparametric convergence rates.

Turning again to our main discussion, note that within our context, the hypotheses of interest are:

\[ H_0 : \max_{j=2, \ldots, m} \int_U E \left( (F_0(u; \theta_0) - F_1(u; \theta_1^j))^2 - (F_0(u; \theta_0) - F_j(u; \theta_j^j))^2 \right) \phi(u) du \leq 0 \]

\[ H_A : \max_{j=2, \ldots, m} \int_U E \left( (F_0(u; \theta_0) - F_1(u; \theta_1^j))^2 - (F_0(u; \theta_0) - F_j(u; \theta_j^j))^2 \right) \phi(u) du > 0. \]

Thus, under \( H_0 \), no model can provide a better approximation than model 1. In order to test \( H_0 \) versus \( H_A \), the relevant test statistic is \( \sqrt{T Z_{T,S}} \), where:

\[ Z_{T,S} = \max_{j=2, \ldots, m} \int_U Z_{j,T,S}(u) \phi(u) du, \tag{1} \]

and

\[ Z_{j,T,S}(u) = \frac{1}{T} \sum_{t=1}^{T} \left( \frac{1}{S} \sum_{n=1}^{S} 1\{Y_{t,n}(\hat{\theta}_{1,T}) \leq u \} \right)^2 \]

\[ - \frac{1}{T} \sum_{t=1}^{T} \left( \frac{1}{S} \sum_{n=1}^{S} 1\{Y_{j,n}(\hat{\theta}_{j,T}) \leq u \} \right)^2, \]

with \( \hat{\theta}_{j,T} \) an estimator of \( \theta_j^j \) that satisfies Assumption 2 below. From equation (1), it is immediate to see that the computational burden of our procedure increases with the dimensionality of \( U \) (i.e. as the number of variables and their lags increases).\(^5\) Unfortunately,

\(^4\)Recently, Giacomini (2005) proposes an extension which uses a weighted (over \( Y_t \)) version of the KLIC, and Kitamura (2002) suggests a generalization for choosing among models that satisfy some conditional moment restrictions.\(^5\) For example, if \( U \) is a two-dimensional subset of \( \mathbb{R}^2 \), and \( \phi \) is uniform on \( U \), then

\[ Z_{T,S} = \frac{1}{N_1 N_2} \max_{j=2, \ldots, m} \sum_{i=1}^{N_1} \sum_{j=1}^{N_2} Z_{j,T,S}(u_{i,j}). \]
Monte Carlo integration techniques, such as importance sampling or one of its accelerated versions (see e.g. Danielsson and Richard (1993)) are not applicable in our context. This is because $Z_{j,T,S}(u)\phi(u)$ is not a joint density and can be either negative or positive. An alternative possibility would be to compute the statistic $\bar{Z}_{T,S} = \max_{j=2,\ldots,m} \frac{1}{T} \sum_{t=1}^{T} \sum_{i=k}^{T} Z_{j,T,S}(Y_{t})$, where $k$ denotes the highest lag order. If $Y_{t}$ were an iid vector-valued process, then $\sqrt{T}Z_{T,S}$ and $\sqrt{T}\bar{Z}_{T,S}$ would be asymptotically equivalent, as shown in Andrews (1997). However, in our case, $Y_{t}$ is a dependent process. Thus, the argument used in Andrews’ proof (based on his Lemma A6) does not apply. Nevertheless, as $T$ gets large, $(Y_{k},Y_{k+1},...Y_{T})$ will become a dense subset in $U$, and so we conjecture that $\sqrt{T}Z_{T,S}$ and $\sqrt{T}\bar{Z}_{T,S}$ are asymptotically equivalent even in the dependent case.

In the sequel, we require the following assumptions:

**Assumption A1**: $Y_{t}$ is stationary-ergodic $\beta-$mixing processes with size $-4$, for $j = 1,\ldots,m$.

**Assumption A2**: For $j = 1,\ldots,m$ : $\sqrt{T} \left( \hat{\theta}_{j,T} - \theta_{j} \right) = A_{j}(\theta_{j}) \frac{1}{\sqrt{T}} \sum_{t=2}^{T} Z_{j,T,S}(Y_{t}) + o_{p}(1)$, where $\frac{1}{\sqrt{T}} \sum_{t=2}^{T} Z_{j,T,S}(Y_{t})$ satisfies a central limit theorem and $A_{j}(\theta_{j})$ is positive definite.\(^6\)

**Assumption A3**: For $j = 1,\ldots,m$ : (i) $\forall \theta_{j} \in \Theta_{j}$, with $\Theta_{j}$ a compact set in $\mathbb{R}^{p_{j}}$ and $Y_{j,n}(\theta_{j})$ is a strictly stationary ergodic $\beta-$mixing process with size $-4$, where $p_{j}$ is the number of estimated parameters in model $j$; (ii) $Y_{j,n}(\theta_{j})$ is continuously differentiable in the interior of $\Theta_{j}$, for $n = 1,\ldots,S$; (iii) $\nabla_{\theta_{j}} Y_{j,n}(\theta_{j})$ is $2r-$dominated in $\Theta_{j}$, uniformly in $n$ for $r > 2$; (iv) $F_{j}(u;\theta_{j})$ is twice continuously differentiable in $u$; and (v) for at least one $j$, $F_{j}(u;\theta_{j}) \neq F_{1}(u;\theta_{1})$ for $u \in \bar{U}$, where $\bar{U}$ is a subset of $U$ of non-zero Lebesgue measure.

**A2** requires that $\sqrt{T} \left( \hat{\theta}_{j,T} - \theta_{j} \right)$ is asymptotically normal with a positive definite covariance matrix. Thus, given the size condition in **A1**, **A2** is satisfied by OLS, NLS, and QMLE, under mild conditions, such as finite $(4+\delta)$th moments and unique identifiability. It is satisfied for the GMM-type estimator of Christiano and Eichenbaum (1992) and the estimator of Bierens (2003). With regard to **A3(i)**, whenever the production function is a Cobb-Douglas type, and the shock to technology follows a unit root process in logs, then output follows

\(^6\)Mixing is a memory requirement stronger than $\alpha-$mixing, but weaker than (uniform) $\phi-$mixing.

\(^7\)Given the size condition in **A1**, **A2** is satisfied by the LS, NLS, QMLE estimator, under mild conditions, such as finite $(4+\delta)$th moments and unique identifiability.

\(^8\)This means that $|\nabla_{\theta_{j}} Y_{j,n}(\theta_{j})| \leq D_{j,n}$, with $\sup_{n} E(D_{j,n}^{2}) < \infty$ (see e.g. Gallant and White (1988), p.33).
a unit root process in logs, and the growth rate is stationary. This is not necessarily true in the case of more general CES production functions. A3(ii) need only hold for estimated parameters. When solving RBC models, we often obtain (linear) ARMA representations for the variables of interest, in terms of final (or reduced form) parameters. Therefore, because of linearity, A3(ii) holds straightforwardly for the final parameters. Hence, if the structural (deep) parameters are smooth functions of the final parameters, as is often the case A3(ii) is satisfied. A3(iii) is a standard assumption (see e.g. Duffie and Singleton (1993)), and A3(iv) is always satisfied for linearized solutions of RBC models. Finally, A3(v) ensures that at least one competing model is nonnested with the benchmark model. This in turn ensures that the covariance matrix of the statistic is positive semi-definite. Hereafter, for notational simplicity, let $F_j(u) = F_j(u; \theta_j^*)$.

**Proposition 1:** Let Assumptions A1-A3 hold. (i) Assume that as $T, S \to \infty : T/S \to \delta$, $0 < \delta < \infty$, then:

$$
\max_{j=2,\ldots,m} \sqrt{T} \int_U \left( Z_{j,T,S}(u) - E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \right) \phi(u) du 
\xrightarrow{d} \max_{j=2,\ldots,m} \int_U Z_j(u) \phi(u) du,
$$

where $Z_j(u)$ is a zero mean Gaussian process with covariance kernal:

$$
K_j(u, u') = 4 \left( F_1(u) - F_j(u) \right) C_{yy}(u, u') \left( F_1(u') - F_j(u') \right) + 4 \mu_{F_1}(u) \delta^2 C_{y_1 y_1}(u, u') \mu_{F_1}(u') \\
+ 4 \mu_{F_2}(u) \delta C_{y_2 y_2}(u, u') \mu_{F_1}(u') + 4 \mu_{F_2}(u) \mu_{f_1, \theta_j}(u) A(\theta_j^*) C_{11} A(\theta_j^*) \mu_{F_1}(u') \mu_{f_1, \theta_j}(u')
+ 4 \mu_{F_1}(u) \mu_{f_1, \theta_j}(u) A(\theta_j^*) C_{1j} A(\theta_j^*) \mu_{F_2}(u') \mu_{f_1, \theta_j}(u') + 8 \mu_{F_1}(u) \delta C_{y_1 y_1}(u, u') (F_1(u') - F_j(u'))
- 8 \mu_{F_2}(u) \delta C_{y_2 y_2}(u, u') (F_1(u') - F_j(u')) + 8 \mu_{F_1}(u) \mu_{f_1, \theta_j}(u) A(\theta_j^*) C_{y_1,1}(u') (F_1(u') - F_j(u'))
- 8 \mu_{F_2}(u) \mu_{f_1, \theta_j}(u) A(\theta_j^*) C_{y_2,1}(u') (F_1(u') - F_j(u')) - 8 \mu_{F_2}(u) \delta C_{y_1 y_1}(u, u') \mu_{F_1}(u')
+ 8 \mu_{F_1}(u) \delta C_{y_1,1}(u') A(\theta_j^*) \mu_{f_1, \theta_j}(u') - 8 \mu_{F_1}(u) \delta C_{y_2,1}(u') A(\theta_j^*) \mu_{F_2}(u') \mu_{f_1, \theta_j}(u')
+ 8 \mu_{F_2}(u) \delta C_{y_2,1}(u') A(\theta_j^*) \mu_{F_2}(u') \mu_{f_1, \theta_j}(u') - 8 \mu_{F_2}(u) \delta C_{y_1,1}(u') A(\theta_j^*) \mu_{F_1}(u') \mu_{f_1, \theta_j}(u')
- 8 \mu_{F_1}(u) \mu_{f_1, \theta_j}(u) A(\theta_j^*) C_{1j} A(\theta_j^*) \mu_{F_2}(u') \mu_{f_1, \theta_j}(u')
$$

with $C_{yy}(u, u') = E \left( \sum_{s=-\infty}^{\infty} \{ Y_2 \leq u \} - F_0(u) \right) \left( \sum_{s=-\infty}^{\infty} \{ Y_{2+s} \leq u \} - F_0(u') \right)$, and where $F_0$ denotes the “true” joint distribution of $Y_t$. Also, $C_{y_j y_j}(u, u') = E \left( \sum_{s=-\infty}^{\infty} \{ Y_{j,2} \leq u \} - F_j(u) \right) \left( \sum_{s=-\infty}^{\infty} \{ Y_{j,2+s} \leq u \} - F_0(u') \right)$. 

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\[ C_{yy}(u, u') = E \left( \sum_{s=-\infty}^{\infty} (1 \{ Y_2 \leq u \} - F_0(u)) (1 \{ Y_{j,2+s} \leq u' \} - F_j(u')) \right), \]
and
\[ C_{jj} = E \left( \sum_{s=-\infty}^{\infty} q_j(Y_2, \theta_j^*) q_j(Y_{2+s}, \theta_j^*) \right), C_{y,j}(u) = E \left( \sum_{s=-\infty}^{\infty} (1 \{ Y_2 \leq u \} - F_j(u)) q_j(Y_{2+s}, \theta_j^*) \right). \]
Finally, \( \mu_{F_j}(u) = E \left( 1 \{ Y_t \leq u \} - F_j(u) \right) \) and \( \mu_{f_i,\theta_j^*}(u) = E \left( f_j(u) \nabla\theta_j Y_{j,n}(\theta_j^*) \right) \), where \( f_j \) is the density of \( F_j \).

(ii) Assume that as \( T, S \to \infty \) \( : S/T^2 \to 0 \) and \( T/S \to 0 \), then:

\[ \max_{j=2, \ldots, m} \sqrt{T} \int_U \left( Z_{j,T,S}(u) - E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \right) \phi(u)du \xrightarrow{d} \max_{j=2, \ldots, m} \int_U \bar{Z}_j(u) \phi(u)du, \]

where \( \bar{Z}_j(u) \) is a zero mean Gaussian process with covariance kernel:

\[ \bar{K}_j(u, u') = 4 (F_1(u) - F_j(u)) C_{yy}(u, u') (F_1(u') - F_j(u')) + 4\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u) A(\theta_j^* \mu_{F_j}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u')) \mu_{f_i,\theta_j^*}(u') \mu_{f_i,\theta_j^*}(u')) + 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u) A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u) A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u) A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u) A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')) - 8\mu_{F_j}(u)\mu_{f_i,\theta_j^*}(u') A(\theta_j^* \mu_{f_i,\theta_j^*}(u') (F_1(u') - F_j(u')). \]

Notice that when \( T/S \to 0 \), then \( \sqrt{T/S} \sum_{n=1}^{S} (1 \{ Y_{j,n}(\theta_j^*) \leq u \} - F_j(u)) \xrightarrow{P} 0 \), uniformly in \( u \), so that the covariance kernel of the limiting distribution does not reflect the contribution of the error term due to the fact we replace the “true” distribution of the simulated series with its empirical counterpart (i.e., simulation error vanishes). From Proposition 1, we see that when all competing models provide an approximation to the true joint distribution that is as accurate as the benchmark, then the limiting distribution is a zero mean Gaussian process with a covariance kernel that reflects the contribution of parameter estimation error, the time series structure of the data and, for \( \delta > 0 \), the contribution of simulation error. This is the case when

\[ \int_U E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \phi(u)du = 0, \text{ for all } j. \]

It follows that in this case, the limiting distribution of

\[ \max_{j=2, \ldots, m} \sqrt{T} \int_U \left( Z_{j,T,S}(u) - E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \right) \phi(u)du \]

is the same as that of \( \sqrt{T} Z_{T,S} \), and so the critical values of the limiting distribution on the RHS of equation (2) provide valid asymptotic critical values for \( \sqrt{T} Z_{T,S} \). On the other hand,
when \( f_U E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \phi(u)du < 0 \) some \( j \), so that at least one alternative model is less accurate than the benchmark, then these critical values can be viewed as upper bounds for critical values from the distribution of \( \sqrt{T}Z_{T,S} \). Note also that when all competing models are less accurate than the benchmark model, the statistic diverges to minus infinity. Finally, under the alternative, \( f_U E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \phi(u)du > 0 \) for some \( j \), so that \( \sqrt{T}Z_{T,S} \) diverges to infinity. Therefore, the test has correct asymptotic size if all models are equally good, is conservative when some model is strictly dominated by the benchmark, and has unit power under the alternative. It should perhaps also be noted, as mentioned in the introduction, that our approach can in principle be modified to allow for the evaluation of predictive densities with rolling and/or recursive estimation schemes (see e.g. Corradi and Swanson (2005b,c,d)).

Although not discussed in the sequel, it should be noted that an “interval” version of the above test also follows directly, given Proposition 1. In particular, if interest focuses on confidence intervals or on testing the null of equal accuracy of two distribution models (analogous to the pairwise conditional mean comparison setup of Diebold and Mariano (1995)), then analogous hypotheses can be constructed by replacing terms of the type \( f_U E \left( (F_0(u; \theta_0) - F_1(u; \theta_1))^2 \right) \phi(u)du \) with terms such as \( E \left( (F_1(\tilde{\omega}; \tilde{\theta}_1) - F_1(\omega; \theta_1)) - (F_0(\tilde{\omega}; \theta_0) - F_0(\omega; \theta_0)) \right)^2 \) or by simply removing \( \max_{j=2,\ldots,m} \) from the statement of the hypotheses, respectively. For example, for the conditional interval version of the above test, we have

\[
H_0 : \max_{j=2,\ldots,m} \mathbb{E} \left( (F_1(\tilde{\omega}; \tilde{\theta}_1) - F_1(\omega; \theta_1)) - (F_0(\tilde{\omega}; \theta_0) - F_0(\omega; \theta_0)) \right)^2
- \left( (F_j(\tilde{\omega}; \tilde{\theta}_j) - F_j(\omega; \theta_j)) - (F_0(\tilde{\omega}; \theta_0) - F_0(\omega; \theta_0)) \right)^2 \leq 0.
\]

versus

\[
H_A : \max_{j=2,\ldots,m} \mathbb{E} \left( (F_1(\tilde{\omega}; \tilde{\theta}_1) - F_1(\omega; \theta_1)) - (F_0(\tilde{\omega}; \theta_0) - F_0(\omega; \theta_0)) \right)^2
- \left( (F_k(\tilde{\omega}; \tilde{\theta}_k) - F_k(\omega; \theta_k)) - (F_0(\tilde{\omega}; \theta_0) - F_0(\omega; \theta_0)) \right)^2 > 0.
\]

Analogous to the statistic outlined above, the appropriate statistic in this case is: \( \sqrt{T}Z_{T,S}^{\text{int}} \), where:

\[
Z_{T,S}^{\text{int}} = \max_{j=2,\ldots,m} \int_{U} Z_{T,S}(u) \phi(u)du,
\]

(3)
and
\[ Z_{j,T,S}^{\text{int}}(u) = \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{ u \leq Y_t \leq \bar{u} \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ u \leq Y_{1,n}(\hat{\theta}_{1,T}) \leq \bar{u} \} \right)^2 - \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{ u \leq Y_t \leq \bar{u} \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ u \leq Y_{j,n}(\hat{\theta}_{j,T}) \leq \bar{u} \} \right)^2. \]

The above test is discussed in Corradi and Swanson (2005a), within the related context of providing a test that allows for the joint comparison of multiple misspecified conditional interval models, for the case of dependent observations. Their test differs from that given above, as conditional distributions are the focus of evaluation. Indeed, there has been growing interest in recent years in providing tests for the correct specification of conditional distributions. One contribution in this direction is the conditional Kolmogorov (CK) test of Andrews (1997), which is based on the comparison of the empirical joint distribution of \( Y_t \) and \( X_t \) with the product of a given distribution of \( y_t | X_t \) and the empirical CDF of \( X_t \) (see also Corradi and Swanson (2005e)). Other contributions in this direction include, for example, Zheng (2000), who suggests a nonparametric test based on a first-order, linear, expansion of the Kullback-Leibler Information Criterion (KLIC), Altissimo and Mele (2002) and Li and Tkacz (2005), who propose a test based on the comparison of a nonparametric kernel estimate of the conditional density with the density implied under the null hypothesis.\(^9\) Following a different route based on use of the probability integral transform, Diebold, Gunther and Tay (1998) suggest a simple and effective means by which predictive densities can be evaluated (see also Bai (2003), Diebold, Hahn and Tay (1999), Hong (2001) and Hong and Li (2005)).

3 Bootstrap Critical Values

In this subsection we outline how to obtain valid critical values for the asymptotic distribution of \( \max_{j=2,\ldots,m} \sqrt{T} \int_U Z_{j,T,S}(u) - E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \phi(u)du \), via use of an empirical process version of the block bootstrap that properly captures the contribution of parameter estimation error, simulation error, and the time series dynamics to the covariance kernel given in Proposition 1. In order to show the first order validity of the bootstrap,

\(^9\)Whang (2000, 2001) proposes a CK type test for the correct specification of the conditional mean.
we derive the limiting distribution of appropriately formed bootstrap statistics and show that they coincide with the limiting distribution given in Proposition 1, recalling that as all candidate models are potentially misspecified under both hypotheses, the parametric bootstrap is not generally applicable in our context. We begin by resampling \( b \) blocks of length \( l \), \( bl = T - 1 \). Let \( Y^*_{t} = (\Delta \log X_t^*, \Delta \log X_{t-1}^*) \) be the resampled series, such that \( Y^*_2, \ldots, Y^*_i, Y^*_i, \ldots, Y^*_{T-1}, \ldots, Y^*_{T} \) equals \( Y_{i,1}, \ldots, Y_{i,1}, Y_{i+1,1}, \ldots, Y_{i+1,1}, \ldots, Y_{i+b,1}, \ldots, Y_{i+b,1} \), where \( I_i, i = 1, \ldots, b \) are independent, discrete uniform random variates on \( 1, \ldots, T - l + 1 \). That is, \( I_i = i, i = 1, \ldots, T - l \) with probability \( 1/(T - l) \). Then, use \( Y^*_{t} \) to compute \( \hat{\theta}^*_{j,T} \) and plug in \( \hat{\theta}^*_{j,T} \) in order to simulate a sample under model \( j, j = 1, \ldots, m \). Let \( Y_{j,n}^*(\hat{\theta}^*_{j,T}), n = 2, \ldots, S \) denote the series simulated in this manner. At this point, we need to distinguish between the case where \( \delta = 0 \) (vanishing simulation error) and \( \delta > 0 \) (non vanishing simulation error). In the former case, we do not need to resample the simulated series, as there is no need to mimic the contribution of simulation error to the covariance kernel. On the other hand, in the latter case we draw \( \tilde{b} \) blocks of length \( \tilde{l} \), with \( \tilde{b}l = S - 1 \), and let \( Y_{j,n}^*(\hat{\theta}^*_{j,T}), j = 1, \ldots, m, n = 2, \ldots, S \) denote the resampled series under model \( j \). Notice that \( Y_{j,n}^*(\hat{\theta}^*_{j,T}), Y_{j,n+1}^*(\hat{\theta}^*_{j,T}), \ldots, Y_{j,S}^*(\hat{\theta}^*_{j,T}) \) is equal to \( Y_{j,\tilde{l}}, \ldots, Y_{j,\tilde{l}+1}(\hat{\theta}^*_{j,T}), \ldots, Y_{j,\tilde{l}+1}(\hat{\theta}^*_{j,T}) \), where \( \tilde{I}_i, i = 1, \ldots, \tilde{b} \) are independent discrete uniform random variates on \( 1, \ldots, S - \tilde{l} \). Also, notice that for each of the \( m \) models, and for each bootstrap replication, we draw \( \tilde{b} \) discrete uniform random variates (the \( \tilde{I}_i \)) on \( 1, \ldots, S - \tilde{l} \), and that draws are independent across models. Thus, in our use of notation, we have suppressed the dependence of \( \tilde{I}_i \) on \( j \).

As discussed above, we consider two different bootstrap versions of \( Z_{T,S}^* \), the first of which is valid when \( T/S \rightarrow \delta > 0 \) and the second of which is valid when \( T/S \rightarrow 0 \), so that \( Y^*_{j,n}^*(\hat{\theta}^*_{j,T}) \) in the first statistic is replaced with \( Y_{j,n}^*(\hat{\theta}^*_{j,T}) \), in the second version. In particular, define \( Z_{T,S}^{*,*} = \max_{j=1,\ldots,m} \int \theta_j \cdot Z_{j,T,S}^*(u) \phi(u) du \), where \( Z_{j,T,S}^*(u) = \frac{1}{T} \sum_{t=1}^{T} \left( \left( 1 \{ Y_t^* \leq u \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ Y_{1,n}^*(\hat{\theta}^*_{1,T}) \leq u \} \right)^2 - \left( 1 \{ Y_t \leq u \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ Y_{1,n}(\hat{\theta}_{1,T}) \leq u \} \right)^2 \right) \)

Also, define \( Z_{T,S}^{*,*} = \max_{j=1,\ldots,m} \int \theta_j \cdot Z_{j,T,S}^*(u) \phi(u) du \), where \( Z_{j,T,S}^*(u) = \frac{1}{T} \sum_{t=1}^{T} \left( \left( 1 \{ Y_t^* \leq u \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ Y_{1,n}^*(\hat{\theta}^*_{1,T}) \leq u \} \right)^2 - \left( 1 \{ Y_t \leq u \} - \frac{1}{S} \sum_{n=1}^{S} 1 \{ Y_{1,n}(\hat{\theta}_{1,T}) \leq u \} \right)^2 \right) \)

Now, let \( \Omega_0 \) be the probability space underlying the historical variables, \( \Omega_j \) the proba-
bility space underlying the series simulated under model $j$ (i.e., the space from which we make random draws in order to form $Y_{j,n}(\hat{\theta}_{j,T})$, where the draws are independent of $\hat{\theta}_{j,T}$, by construction). Then, define the enlarged probability space, $\Omega$, as $\Omega = \Omega_0 \times \Omega_1 \times \ldots \times \Omega_m$. Also, let $P_0$ be the probability law governing the actual sample, and $P_j$ be the probability law of the samples simulated under model $j$. That is, $P_j$ is the probability law governing $Y_{j,n}(\theta_j)$. Now, define $P = P_0 \times P_1 \times \ldots \times P_m$. Analogously, define $P_0^*$ to be the probability law governing the resampled series, $Y_t^*$, conditional on the historical sample, $Y_t$, and define $P_j^*$ to be the probability law governing the resampled series, $Y_{j,n^*}(\theta_j)$, conditional on the simulated series, $Y_{j,n}(\theta_j)$. Now, define $P^* = P_0^* \times P_1^* \times \ldots \times P_m^*$. Finally, let $\omega$ denote a draw from $\Omega$, and let $I_T$ and $I_S$ denote the block length of the resampled series from the actual and simulated samples, respectively. Note that when constructing $Z_{j,T,S}^*$, $\delta = 0$ so that $P^* = P_0$.

**Proposition 2:** Let Assumptions A1-A3 hold. Also, assume that as $I_T, I_S \to \infty : I_T/T^{1/2} \to 0$ and $I_S/S^{1/2} \to 0$. (i) Assume that as $T, S \to \infty : T/S \to \delta$, $0 < \delta < \infty$, then:

$$P\left(\omega : \sup_{v \in \mathbb{R}} \left| P^* \left( \max_{j=2,\ldots,m} \sqrt{T} \int_U Z_{j,T,S}^*(u) \phi(u) du \leq v \right) \right| > \varepsilon \right) \to 0.$$ 

(ii) Assume that as $T, S \to \infty : S/T^2 \to 0$ and $T/S \to 0$, then:

$$P\left(\omega : \sup_{v \in \mathbb{R}} \left| P^* \left( \max_{j=2,\ldots,m} \sqrt{T} \int_U Z_{j,T,S}^*(u) \phi(u) du \leq v \right) \right| > \varepsilon \right) \to 0.$$

Thus, for any bootstrap replication, compute the bootstrap statistic, $\sqrt{T}Z_{j,T,S}^*$ ($\sqrt{T}Z_{T,S}^*$). Perform $B$ bootstrap replications ($B$ large) and compute the quantiles of the resultant empirical distribution. Reject $H_0$ if $\sqrt{T}Z_{T,S}$ is greater than the $(1-\alpha)th$-quantile. Otherwise, do not reject. Now, for all samples except a set with probability measure approaching zero, $\sqrt{T}Z_{T,S}$ has the same limiting distribution as the corresponding bootstrapped statistic, when $\int_U E \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right) \phi(u) du = 0$ for all $j = 2,\ldots,m$. In this case, the above approach ensures that the test has asymptotic size equal to $\alpha$. On the other hand,

---

10Note that although we generate the simulated series using $\hat{\theta}_{j,T}$, $P_j$ denotes the probability law of the series simulated under $\theta_j$, which is the probability limit of $\hat{\theta}_{j,T}$.
when one (or more) competing models is (are) strictly dominated by the benchmark, the approach ensures that the test has an asymptotic size between 0 and $\alpha$. Finally, under the alternative, $Z_{T,S}$ diverges to (plus) infinity, while the corresponding bootstrap statistic has a well defined limiting distribution. This ensures unit asymptotic power. As our bootstrap mimics the limiting distribution of the statistics in the least favorable case for the null, inference is conservative (see Hansen (2005) for further discussion).

4 Empirical Illustration

Consider a stochastic growth model characterized by flexible labor supply, non-zero capital depreciation, and two shocks. The first is the usual permanent shock to technology, and the second is a transitory preference (or taste) shock. The model is exactly that considered by Christiano (1988) and the accompanying appendix (i.e. Christiano (1987)), except that we use a Cobb-Douglas production function (i.e. we set $\sigma = 0$ in his equation (2)), so that inventories do not play a role in production. Assume that the representative agent chooses a contingent plan for consumption and labor supply which maximizes

$$\mathbb{E}_0 \sum_{t=0}^{\infty} \beta^t (\exp(u_t) \log C_t - \gamma H_t),$$

where $C_t$ and $H_t$ are per capita consumption and hours worked, respectively. Additionally, $u_t$ is a transitory taste shock defined in (7) below. The maximization above is subject to the production constraint,

$$C_t \leq X_t - K_t + \frac{1-\delta}{n} K_{t-1},$$

where $X_t$ is per capita output, $K_t$ per capita capital stock, $n$ is the rate of growth of population (assumed constant), $\delta$ is the depreciation rate. Additionally,

$$X_t = (z_t H_t)^{1-\theta} K_{t-1}^\theta,$$

where $z_t$ is a permanent technological shock, defined as $z_t = z_{t-1} \exp(t_t)$, and where $\theta$ denotes the capital share. Now, assume that

$$
\begin{pmatrix}
  u_t \\
  t_t
\end{pmatrix} =
\begin{pmatrix}
  0 & 0 \\
  a_2 & A_{22}
\end{pmatrix}
\begin{pmatrix}
  u_{t-1} \\
  t_{t-1}
\end{pmatrix} +
\begin{pmatrix}
  \epsilon_{1t} \\
  \epsilon_{2t}
\end{pmatrix},
\end{equation}$$
where $-1 < A_{11}, A_{22} < 1$. The structural parameters in this model are: $\beta, \delta, n, \theta, \gamma, a_2, A_{11},$ and $A_{22}$. As we cannot find explicit solution of (4) subject to (5), we instead solve the following quadratic approximation:

$$\max_{d_t} E_0 \sum_{t=0}^{\infty} \beta^t (c_1' d_t + c_2' s_t + s'_t R s_t + d'_t Q d_t + 2s'_t F d_t)$$

subject to $s_{t+1} = \phi_0 + \phi_1 s_t + B d_t + \epsilon_t$, where $s_t^* = (k_{t-1}, u_t, t_t, u_{t-1}, t_{t-1})'$ and $d_t^* = (k_t^*, h_t^*)'$ with $k_t^* = \log(K_t/z_{t-1})$ and $h_t^* = \log(H_t)$; and where $c_1, c_2, R, Q, F, \phi_0, \phi_1, B$ are parameters constructed using functions of first and second derivatives. (For complete details, see Appendix B of Corradi and Swanson (2003).) Assume that our objective is to find the dynamics of $X_t$ and $H_t$, as given by the constrained maximization problem above. The optimal policy function governing the dynamics of $d_t^*$ are given by,

$$d_t^* = K_0 + K_1 s_t^*,$$

where $K_0 = -\frac{1}{2}Q^{-1}c_1$, $K_1 = -Q^{-1}F'$, $K_0$ is a $2 \times 1$ vector and $K_1$ is a $2 \times 5$ matrix. In this framework, it follows from Christiano (1988) that

$$\tau'_1 = c'_1 + \beta V'_s B + 2\beta \phi'_0 V_Q B,$$

$$Q = Q + \beta B'V_Q B,$$ and

$$F = F + \beta \phi'_1 V_Q B.$$ (12)

In order to determine (9), we need to construct $K_0$ and $K_1$ by solving 2 Riccati equations for $V_Q$ and $V_s$ based on (10), (11), and (12). We begin with the determination of $V_Q$. $V_Q$ is the solution of the following Riccati equation:

$$V_{Qt} = R + \beta \phi'_1 V_{Qt+1} \phi_1 - (\beta \phi'_1 V_{Qt+1} B + F) \times (Q + \beta B'V_{Qt+1} B)^{-1} (\beta B'V_{Qt+1} \phi_1 + F')$$

(13)

An explicit closed-form solution to this Riccati equation does not exist. However we can find an arbitrarily accurate solution via the iteration method discussed in MacGrattan (1990). In
particular, starting with an initial value, say $V_{Q,0}$ (typically set equal to the identity matrix time a small number), at the $n + 1$-th iteration, we have

$$V_{Q}^{n+1} = R + \beta \phi_1 V_{Q}^{n} \phi_1 - (\beta \phi_1 V_{Q}^{n} B + F) \times \left( Q + \beta B' V_{Q}^{n} B \right)^{-1} \left( \beta B' V_{Q}^{n} \phi_1 + F' \right).$$

The stopping point occurs at that value of $V_{Q}^{n+1}$ for which $\| V_{Q}^{n+1} - V_{Q}^{n} \| < \varepsilon$, where $\| \cdot \|$ denotes the euclidean norm, and where we set $\varepsilon = 0.0000001$. Now, in much the same manner, and by using $V_{Q}^{n+1}$ we can obtain an (approximate) solution for the Riccati equation describing $V_s$. Namely,

$$V_{s}^{n+1} = (I - \beta (K_1 B' + \phi_1'))^{-1} \left( K_1 \left( c_1 + 2 \beta B' V_{Q}^{n+1} \phi_0 \right) + c_2 + 2 \beta \phi_1 V_{Q}^{n+1} \phi_0 \right).$$

Now,

$$d_t^* = \begin{pmatrix} k_{t}^* \\ h_t^* \end{pmatrix} = K_0 + K_1 s_t^* = \begin{pmatrix} k_{01} \\ k_{02} \end{pmatrix} + \begin{pmatrix} k_{11} & k_{12} & k_{13} & k_{14} & k_{15} \\ k_{21} & k_{22} & k_{23} & k_{24} & k_{25} \end{pmatrix} \begin{pmatrix} k_{t-1}^* \\ u_t \\ t_t \\ u_{t-1} \\ t_{t-1} \end{pmatrix},$$

where $K_0$ and $K_1$ are defined as above. Thus,

$$\begin{pmatrix} k_{t}^* \\ h_t^* \end{pmatrix} = \begin{pmatrix} k_{01} + k_{13} a_2 \\ k_{02} + k_{23} a_2 \end{pmatrix} + \begin{pmatrix} k_{11} \\ k_{21} \end{pmatrix} k_{t-1}^* + \begin{pmatrix} k_{12} A_{11} + k_{14} & k_{13} A_{22} + k_{15} \\ k_{22} A_{11} + k_{24} & k_{23} A_{22} + k_{25} \end{pmatrix} \begin{pmatrix} u_{t-1} \\ t_{t-1} \end{pmatrix}$$

$$+ \begin{pmatrix} k_{12} & k_{13} \\ k_{22} & k_{23} \end{pmatrix} \begin{pmatrix} \epsilon_{1t} \\ \epsilon_{2t} \end{pmatrix}$$

Finally, recalling that $k_t^* = \log(K_t/z_{t-1})$ and $h_t^* = \log(H_t)$, and recalling (7), we can simulate capital and hours using:

$$\begin{pmatrix} \log K_t \\ \log H_t \end{pmatrix} = \begin{pmatrix} \log z_{t-1} \\ 0 \end{pmatrix} + \begin{pmatrix} k_{01} + k_{13} a_2 \\ k_{02} + k_{23} a_2 \end{pmatrix} + \begin{pmatrix} k_{11} \\ k_{21} \end{pmatrix} \left( \log K_{t-1} - \log z_{t-2} \right)$$

$$+ \begin{pmatrix} k_{12} A_{11} + k_{14} & k_{13} A_{22} + k_{15} \\ k_{22} A_{11} + k_{24} & k_{23} A_{22} + k_{25} \end{pmatrix} \begin{pmatrix} u_{t-1} \\ t_{t-1} \end{pmatrix} + \begin{pmatrix} k_{12} & k_{13} \\ k_{22} & k_{23} \end{pmatrix} \begin{pmatrix} \epsilon_{1t} \\ \epsilon_{2t} \end{pmatrix}$$

(18)
Note that all of the parameters in (18) are explicit functions of the structural parameters \( \beta, \delta, n, \theta, \gamma, a_2, A_{11}, \) and \( A_{22}. \) Note also that for \( |k_{11}| < 1, \log K_t - \log z_{t-1} \) is \( I(0), \) and therefore, for \( |k_{21}| < 1, \log H_t \) is \( I(0) \) too. Also, as \( \log z_{t-1} \) is \( I(1), \log K_t \) is \( I(1), \) so that both capital and output (see (19) below) follow unit root processes in logs. Finally, from the production function in (6), upon taking logs,

\[
\log X_t = (1 - \theta) \log z_t - \theta \log n + (1 - \theta) \log H_t + \theta \log K_{t-1} \tag{19}
\]

In the sequel, we take values for structural parameters \( \beta, \delta, n, \theta, \gamma, a_2, A_{11}, \) and \( A_{22} \) from King, Plosser and Rebelo (1988a,b) and Christiano (1988). In particular, and unless otherwise stated, we set \( \beta = 0.99, n = 1.012, \delta = 0.025, \theta = 0.50, \gamma = 0.0028, a_2 = 0.0044, A_{22} = -0.10, A_{11} = 0.95. \) Furthermore, and unless otherwise stated, we estimate \( \sigma^2_1 \) and \( \sigma^2_2 \) by simply constructing simulations with various values for these variances and minimizing the difference between the simulated and actual variance of per capita output and hours growth.\(^{11}\)\(^{12}\) Given this setup, we evaluate the joint distribution of \( Y_t = (\log \Delta X_t, \log H_t) \) as well as \( Y_i = (\log \Delta X_t, \log H_t, \log \Delta X_{t-1}, \log H_{t-1}). \)

Some illustrative results are given in Tables 1 and 2. The first column of numerical entries reports \( \sqrt{T}Z_{T,S} \) for \( T = 70 \) (corresponding to the 70 annual U.S. per capita output growth observations used for the period 1933-2002), and for \( S = \{T, 2T, 5T, 10T\}. \)\(^{13}\) All statistics are based on use of 20 uniformly distributed values for \( u \) across the range of the actual data. The next four columns of entries contain 5\% and 10\% bootstrap critical values constructed using both \( \sqrt{T}Z_{T,S}^{**} \) and \( \sqrt{T}Z_{T,S}^* \) (i.e. \( T/S - > \delta > 0 \) and \( T/S- > 0 \), respectively). For the case where \( S = T, \) we set the bootstrap block length equal to \( l_1 = 2, l_2 = 5, \) and \( l_3 = 7. \) For all other cases, where \( S = aT, \) say, we set \( l_i \) equal to “\( a \)” times the corresponding value of \( l_i \) when \( S = T, \) for \( i = 1, 2, 3. \) Thus, “\( a \)” takes the values 2.5, and 10, depending on

\(^{11}\) Throughout, and for simplicity, we assume that the covariance between the shocks is zero. This assumption is not inconsistent with some of the empirical findings in the literature.

\(^{12}\) Note that as we are interested in matching the covariance structure between historical and simulated data, we can use different specifications of the marginal distribution of the shocks. In the current context, the variance estimators we use depend also the simulated sample size \( S. \) Call these estimators \( \hat{\xi}_{T,S}. \) If \( S/T \rightarrow 0, \) then simulation error is negligible and Assumption A2 is satisfied as \( \hat{\xi}_{T,S} \) is asymptotically equivalent to GMM. In the case where \( S/T \rightarrow 0, \) then we also require that the scaled difference between the simulated and actual sample moments is asymptotically normal. This is in general the case, provided that Assumption A3 holds.

\(^{13}\) Note that in all case \( S/T > 0, \) however, when for example \( T = 70 \) and \( S = 700, \) then simulation error is likely to be negligible with respect to parameter estimation error.
the magnitude of $S$. Bootstrap empirical distributions are constructed using 100 bootstrap replications (i.e. $B = 100$). Even though hours and output are examined, for the sake of brevity we only discuss output in the sequel. Various summary statistics (including $\% < 0$, min, max, and standard error) constructed using the simulated data are given in the remaining columns of the tables. Corresponding values based on the actual data are: $\% < 0 = 0.213$; min = -0.128; max = 0.158; mean = 0.027; and standard error = 0.044. Of course, the $Z_{T,S}$ statistics are based on the joint evaluation of all variables in $Y_t$, including output and hours. We consider four different scenarios: Part I - Compare 3 models using $Y_t = (\log \Delta X_t, \log H_t)$. Set $\beta = \{0.95, 0.90, 0.99\}$. All shocks are normally distributed with variance estimated as discussed above. The model with $\beta = 0.99$ is the benchmark. Part II - Compare 3 models using $Y_t = (\log \Delta X_t, \log H_t)$. The benchmark has shocks drawn from a $1/3\chi_3$ distribution. One alternative assumes normality and uses estimated variances while the other assumes normality and sets $\sigma_{\epsilon_1}^2 = 0.039$ and $\sigma_{\epsilon_2}^2 = 0.077$ (as in Christiano (1988)). Part III - As Part I, except that $Y_t = (\Delta \log X_t, \log H_t, \Delta \log X_{t-1}, \log H_{t-1})$. Part IV - As Part II, except that $Y_t = (\Delta \log X_t, \log H_t, \Delta \log X_{t-1}, \log H_{t-1})$.

Turning to our results, notice in Part I of Table 1 that the null hypothesis fails to be rejected in almost all cases, and never rejects when the simulation sample is 5 or 10 times as large as the sample size. This suggests that the discount rate can equally take values between 0.90 and 0.99, at least when the objective is to match the joint dynamics of simulated output and hours as closely as possible with the historical record. However, notice in Part II of Table 1 that the benchmark where shocks follow Student’s t-distributions is soundly rejected in every case. Furthermore, the dynamics of output are clearly far from reasonable under the benchmark as well as under the first alternative model (see Sim1 and Sim2 summary statistics), corresponding to the t-distribution assumption (Sim1) and the case where we assume normality and fix $\sigma_{\epsilon_1}^2 = 0.039$ and $\sigma_{\epsilon_2}^2 = 0.077$ (Sim2). This supports the use of our simple simulated GMM type scheme for estimating the shock variances (see Sim3 results). Table 2 contains results for experiments analogous to those reported in Table 1, except that the joint dynamics of four variables are examined. The results of these experiments are the same as those reported on in Table 1, suggesting that generalizing our approach to the assessment of more variables does not affect test performance.
5 Concluding Remarks

In this paper we propose a test for comparing the joint distributions of historical time series with those simulated under a given DSGE model via a distributional generalization of the reality check of White (2000) in which we assess whether competing models are more accurate approximations to the “true” distribution than a given benchmark model, in a squared error sense. Two empirical versions of the block bootstrap are used to construct valid asymptotic critical values. Finally, an illustrative example is given in which the testing approach is applied to an RBC model, and it is found that RBC models are quite sensitive to distributional and error variance magnitude assumptions, but are less sensitive to small changes in primitive parameter values.
6 Appendix

Hereafter, $P$, $P^*$, $\Omega$, $P_i$, $P_i^*$ and $\Omega_i$, $i = 0, 1, \ldots, m$ are defined above the statement of Proposition 2. Also, $o_p(1)$ and $o_{P_i^*}(1)$ denote terms approaching zero in probability $P^*$, $\forall \omega \in \Omega$ and $P_i^*$, $\forall \omega_i \in \Omega_i$, respectively. Furthermore, $E_{P^*}$ and $\text{Var}_{P_i^*}$ are mean and variance operators with respect to $P^*$, with $E_{P_i^*}$ and $\text{Var}_{P_i^*}$ defined analogously. Finally, $csz$ is meant to mean “conditional on the sample and for all samples except a set of zero probability measure”, while $cszp$ is the same, except applies to all samples, simulated and empirical.

Proof of Proposition 1: (i) We begin by examining the limiting distribution, pointwise in $u$. Now,

$$\sqrt{T} Z_{j,T,s}(u) = \frac{1}{\sqrt{T}} \sum_{t=1}^{T} \left( (1 \{ Y_t \leq u \} - F_1(u)) - \left( \frac{1}{S} \sum_{n=1}^{S} 1 \{ Y_{1,n}^*(\theta^*_1) \leq u - (Y_{1,n}^*(\hat{\theta}_{1,T}) - Y_{1,n}^*(\theta^*_1)) \} \right) \right.$$ 

$$- F_1(u - (Y_{1,n}^*(\hat{\theta}_{1,T}) - Y_{1,n}^*(\theta^*_1))) \right) \right)^2 - \frac{1}{\sqrt{T}} \sum_{t=1}^{T} \left( (1 \{ Y_t \leq u \} - F_j(u)) - \frac{1}{S} \sum_{n=1}^{S} \left( 1 \{ Y_{j,n}^*(\theta^*_j) \leq u - (Y_{j,n}^*(\hat{\theta}_{j,T}) - Y_{j,n}^*(\theta^*_j)) \} \right) \right.$$

$$- F_j(u - (Y_{j,n}^*(\hat{\theta}_{j,T}) - Y_{j,n}^*(\theta^*_j))) \right) \right)^2 \right)^2$$

Given A3(i), $\frac{1}{\sqrt{T}} \sum_{n=1}^{S} (1 \{ Y_{j,n}^*(\theta^*_j) \leq u \} - F_j(u))$ is stochastic equicontinuous in $u \in U$, from Theorem 1 (Application 1) of Doukhan, Massart and Rio (1995). Also, given A2 and A3(ii),

$$Y_{j,n}^*(\hat{\theta}_{j,T}) - Y_{j,n}^*(\theta^*_j) = \nabla_{\theta_j} Y_{j,n}^*(\hat{\theta}_{j,T}) (\hat{\theta}_{j,T} - \theta^*_j) = \frac{1}{\sqrt{T}} \nabla_{\theta_j} Y_{j,n}^*(\theta_{j,T}) O_P(1), \quad 14$$

with $\theta_{j,T} \in (\hat{\theta}_{j,T}, \theta^*_j)$. Now, given the domination conditions in A3(iii), by Chebyshev’s inequality,

$$\text{Pr} \left( \frac{1}{\sqrt{T}} \sup_{n \leq S} | \nabla_{\theta_j} Y_{j,n}^*(\theta_j) | > \varepsilon \right) \leq \sum_{i=1}^{S} \text{Pr} \left( \frac{1}{\sqrt{T}} | \nabla_{\theta_j} Y_{j,n}^*(\theta_j) | > \varepsilon \right)$$

$$\leq \sum_{i=1}^{S} \text{Pr} \left( \frac{1}{\sqrt{T}} D_j > \varepsilon \right) \leq \frac{1}{\varepsilon^2} \text{Pr} \left( E(D_j^2) \rightarrow 0 \right)$$
as \( S/T^2 \to 0 \), given that \( S/T \to \delta > 0 \). Thus,

\[
\sqrt{\frac{T}{S}} \frac{1}{\sqrt{S}} \sum_{n=1}^{S} \left( 1 \{Y_{1,n}(\theta^*_1) \leq u - (Y_{1,n}(\widehat{\theta}_{1,T}) - Y_{1,n}(\theta^*_1)) \} - \left( F_1(u - (Y_{1,n}(\widehat{\theta}_{1,T}) - Y_{1,n}(\theta^*_1))) - F_1(u) \right) \right) = \sqrt{\frac{T}{S}} \frac{1}{\sqrt{S}} \sum_{n=1}^{S} \left( 1 \{Y_{1,n}(\theta^*_1) \leq u \} - F_1(u) \right) + o_P(1),
\]

with the \( o_P(1) \) term holding uniformly in \( u \). Noticing that \( F_j(u - (Y_{j,n}(\widehat{\theta}_{j,T}) - Y_{j,n}(\theta^*_j))) = f_j(u - (Y_{j,n}(\widehat{\theta}_{j,T}) - Y_{j,n}(\theta^*_j))) \nabla_{\widehat{\theta}_j} Y_{j,n}(\widehat{\theta}_{j,T})' \left( \widehat{\theta}_{j,T} - \theta^*_j \right) \), where \( f_j \) is the density associated with \( F_j \), after some elementary manipulations, the right hand side (RHS) of (20) can be written as:

\[
- \frac{1}{S} \sum_{n=1}^{S} f_1(u - (Y_{1,n}(\widehat{\theta}_{1,T}) - Y_{1,n}(\theta^*_1))) \nabla_{\widehat{\theta}_1} Y_{1,n}(\widehat{\theta}_{1,T})' \left( \widehat{\theta}_{1,T} - \theta^*_1 \right) \right)^2
\]

\[
- \frac{1}{\sqrt{T}} \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{Y_t \leq u \} - F_0(u) \right) - (F_j(u) - F_0(u)) - \left( \frac{1}{S} \sum_{n=1}^{S} 1 \{Y_{1,n}(\theta^*_1) \leq u \} - F_1(u) \right)
\]

\[
- \frac{1}{S} \sum_{n=1}^{S} f_j(u - (Y_{j,n}(\widehat{\theta}_{j,T}) - Y_{j,n}(\theta^*_j))) \nabla_{\widehat{\theta}_j} Y_{j,n}(\widehat{\theta}_{j,T})' \left( \widehat{\theta}_{j,T} - \theta^*_j \right) \right)^2 + o_P(1),
\]

where the \( o_P(1) \) term holds uniformly in \( u \). Now,

\[
\sqrt{T} \left( \frac{1}{S} \sum_{n=1}^{S} f_j(u - (Y_{j,n}(\widehat{\theta}_{j,T}) - Y_{j,n}(\theta^*_j)))^2 \nabla_{\widehat{\theta}_j} Y_{j,n}(\widehat{\theta}_{j,T})' \left( \widehat{\theta}_{j,T} - \theta^*_j \right) \right)^2 = o_P(1),
\]

(21)

uniformly in \( u \), as \( (\widehat{\theta}_{j,T} - \theta^*_j) = O_P(T^{-1/2}) \). Given A1 and A3, uniformly in \( u \),

\[
\sqrt{T} \left( \frac{1}{S} \sum_{n=1}^{S} 1 \{Y_{j,n}(\theta^*_j) \leq u \} - F_j(u) \right)^2 = O_P \left( \frac{\sqrt{T}}{S} \right) = o_P(1),
\]

(22)

and

\[
\frac{1}{S} \sum_{n=1}^{S} \left( 1 \{Y_{j,n}(\theta^*_j) \leq u \} - F_j(u) \right) \frac{1}{S} \sum_{n=1}^{S} f_j(u - (Y_{j,n}(\widehat{\theta}_{j,T}) - Y_{j,n}(\theta^*_j))) \nabla_{\widehat{\theta}_j} Y_{j,n}(\widehat{\theta}_{j,T})' \sqrt{T} \left( \widehat{\theta}_{j,T} - \theta^*_j \right) \]

\[
= o_P(1).
\]

(23)
Therefore, pointwise in $K$ (ii) continuous mapping theorem. Application 1 in Doukh, Massart and Del Rio (1995)). The desired result follows by the Cramer-Wold device. Finally, given A1 and A2, by the uniform law of large numbers for mixing processes,

$$
\sqrt{T}Z_{j,T,S}(u) - \sqrt{T} \left( (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right)
$$

$$
= -\frac{2}{\sqrt{T}} \sum_{t=1}^{T} (1 \{ Y_t \leq u \} - F_0(u)) (F_1(u) - F_j(u))
$$

$$
-2\mu_{F_1}(u)\delta_1 \frac{1}{S} \sum_{s=1}^{S} \left( 1 \{ Y_{1,n}(\theta_1^s) \leq u \} - F_1(u) \right) + 2\mu_{F_j}(u)\delta_1 \frac{1}{S} \sum_{s=1}^{S} \left( 1 \{ Y_{j,n}(\theta_j^s) \leq u \} - F_j(u) \right)
$$

$$
-2\mu_{F_1}(u)\mu_{F_j}(u)\delta_1 \frac{1}{S} \sqrt{T} \left( \hat{\theta}_1,T - \theta_1^s \right) + 2\mu_{F_j}(u)\mu_{F_j}(u)\delta_1 \frac{1}{S} \sqrt{T} \left( \hat{\theta}_j,T - \theta_j^s \right) + o_p(1).
$$

Therefore, pointwise in $u$, $\sqrt{T} \left( Z_{j,T,S}(u) - (F_0(u) - F_1(u))^2 - (F_0(u) - F_j(u))^2 \right)$ satisfies a CLT, and the asymptotic variance is as given in the statement of the proposition. Convergence of the finite dimensional distribution follows immediately upon application of the Cramer-Wold device. Finally, given A1 and A3, $\frac{2}{\sqrt{T}} \sum_{t=1}^{T} (1 \{ Y_t \leq u \} - F_0(u))$ and $\frac{1}{\sqrt{T}} \sum_{t=1}^{T} (1 \{ Y_{j,n}(\theta_1^s) \leq u \} - F_j(u))$ are stochastic equicontinuous in $u$ (e.g. by Theorem 1, Application 1 in Doukh, Massart and Del Rio (1995)). The desired result follows by the continuous mapping theorem.

(ii) The proof follows using the same argument as in part (i), and when the expression for $K_j(u,u')$ in part (i) is adjusted by setting $\delta = 0$.

**Proof of Proposition 2:** (i) Note that:

$$
\sqrt{T}Z_{j,T,S}(u) = \frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( \left( 1 \{ Y_t \leq u \} - F_1(u) \right) - \frac{1}{S} \sum_{n=1}^{S} \left( 1 \{ Y_{1,n}^*(\hat{\theta}_1,T) \leq u \} - F_1(u) \right) \right)^2
$$

$$
\left( 1 \{ Y_t \leq u \} - F_1(u) \right) - \frac{1}{S} \sum_{n=1}^{S} \left( 1 \{ Y_{1,n}(\hat{\theta}_1,T) \leq u \} - F_1(u) \right)
$$

$$
\frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( \left( 1 \{ Y_t \leq u \} - F_j(u) \right) - \frac{1}{S} \sum_{n=1}^{S} \left( 1 \{ Y_{j,n}^*(\hat{\theta}_j,T) \leq u \} - F_j(u) \right) \right)^2
$$

$$
\left( 1 \{ Y_t \leq u \} - F_j(u) \right) - \frac{1}{S} \sum_{n=1}^{S} \left( 1 \{ Y_{j,n}(\hat{\theta}_j,T) \leq u \} - F_j(u) \right)
$$

$$
. \quad (24)
$$
Now, we begin by considering the first term on the RHS of (24), which can be rewritten as,

\[
\begin{align*}
= \frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( 1\{Y_t^* \leq u\} - F_1(u) \right) &- \frac{1}{S} \sum_{n=1}^{S} \left( 1\{Y_{1,n}^* (\theta^*_1) \leq u - (Y_{1,n}^* (\hat{\theta}^*_{1,T}) - Y_{1,n}^* (\theta^*_1)) \} \right) \\
&- F_1(u) - (Y_{1,n}^* (\hat{\theta}^*_{1,T}) - Y_{1,n}^* (\theta^*_1))) - \frac{1}{S} \sum_{n=1}^{S} \left( F_1(u) - (Y_{1,n}^* (\hat{\theta}^*_{1,T}) - Y_{1,n}^* (\theta^*_1))) - F_1(u) \right)^2 \\
&- \frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( 1\{Y_t \leq u\} - F_1(u) \right) - \frac{1}{S} \sum_{n=1}^{S} \left( 1\{Y_{1,n} (\theta^*_1) \leq u - (Y_{1,n} (\hat{\theta}^*_{1,T}) - Y_{1,n} (\theta^*_1))) \} \right) \\
&- F_1(u) - (Y_{1,n} (\hat{\theta}^*_{1,T}) - Y_{1,n} (\theta^*_1))) - \frac{1}{S} \sum_{n=1}^{S} \left( F_1(u) - (Y_{1,n} (\hat{\theta}^*_{1,T}) - Y_{1,n} (\theta^*_1))) - F_1(u) \right)^2 \\
= \frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( 1\{Y_t^* \leq u\} - F_1(u) \right) &- \frac{1}{S} \sum_{n=1}^{S} \left( 1\{Y_{1,n}^* (\theta^*_1) \leq u \} - F_1(u) \right) \\
&- \frac{1}{S} \sum_{n=1}^{S} f_1 \left( u - (Y_{1,n}^* (\bar{\theta}^*_{1,T}) - Y_{1,n}^* (\theta^*_1))) \right) \nabla_{\theta^*_1} Y_{1,n}^* (\bar{\theta}^*_{1,T}) \sqrt{T} \left( \hat{\theta}^*_{1,T} - \theta^*_1 \right)^2 \\
&- \frac{1}{T^{1/2}} \sum_{t=1}^{T} \left( 1\{Y_t \leq u\} - F_1(u) \right) &- \frac{1}{S} \sum_{n=1}^{S} \left( 1\{Y_{1,n} (\theta^*_1) \leq u \} - F_1(u) \right) \\
&- \frac{1}{S} \sum_{n=1}^{S} f_1 \left( u - (Y_{1,n} (\bar{\theta}^*_{1,T}) - Y_{1,n} (\theta^*_1))) \right) \nabla_{\theta^*_1} Y_{1,n} (\bar{\theta}^*_{1,T}) \sqrt{T} \left( \hat{\theta}^*_{1,T} - \theta^*_1 \right)^2 \\
&+ o_P(1), \quad \Pr - P.
\end{align*}
\]

The second term on the RHS of (24) can be treated in the same way, by replacing 1 with j.

Therefore, after some elementary manipulations, \(Z_{j,T,S}^*(u)\) can be written as:

\[
Z_{j,T,S}^*(u) = - \frac{2}{\sqrt{T}} \sum_{t=1}^{T} \left( \{Y_t^* \leq u\} - \{Y_t \leq u\} \right) \left( F_1(u) - F_j(u) \right) \\
- d_{1,1} \sqrt{\frac{T}{S}} \sum_{n=1}^{S} \left( 1\{Y_{1,n}^*(\theta^*_1) \leq u\} - F_1(u) \right) + d_{2,1} \sqrt{\frac{T}{S}} \sum_{n=1}^{S} \left( 1\{Y_{1,n}(\theta^*_1) \leq u\} - F_1(u) \right)
\]
\[ +d_{1,j} \sqrt{\frac{T}{S}} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} \left( 1 \{ Y_{j,n}^* (\theta_j^*) \leq u \} - F_j(u) \right) - d_{2,j} \sqrt{\frac{T}{S}} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} \left( 1 \{ Y_{j,n} (\theta_j^*) \leq u \} - F_j(u) \right) \]

\[ -d_{1,1} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} f_1 \left( u - (Y_{1,n}^*(\overline{\theta}_1) - Y_{1,n}(\theta_1^*)) \right) \nabla_{\theta_1} Y_{1,n}^*(\overline{\theta}_1) \sqrt{T} \left( \hat{\theta}_1^* - \theta_1^* \right) \]

\[ +d_{2,1} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} f_1 \left( u - (Y_{1,n}(\overline{\theta}_1) - Y_{1,n}(\theta_1^*)) \right) \nabla_{\theta_1} Y_{1,n}(\overline{\theta}_1) \sqrt{T} \left( \hat{\theta}_1 - \theta_1^* \right) \]

\[ +d_{1,j} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} f_j \left( u - (Y_{j,n}^*(\overline{\theta}_j) - Y_{j,n}(\theta_j^*)) \right) \nabla_{\theta_j} Y_{j,n}^*(\overline{\theta}_j) \sqrt{T} \left( \hat{\theta}_j^* - \theta_j^* \right) \]

\[ -d_{2,j} \frac{1}{\sqrt{s}} \sum_{n=1}^{s} f_j \left( u - (Y_{j,n}(\overline{\theta}_j) - Y_{j,n}(\theta_j^*)) \right) \nabla_{\theta_j} Y_{j,n}(\overline{\theta}_j) \sqrt{T} \left( \hat{\theta}_j - \theta_j^* \right) \]

where \( d_{1,x} = \frac{2}{T} \sum_{t=1}^{T} \left( 1 \{ Y_t^* \leq u \} - F_x(u) \right) \) and \( d_{2,x} = \frac{2}{T} \sum_{t=1}^{T} \left( 1 \{ Y_t \leq u \} - F_x(u) \right) \), \( x = 1, j \). Now, conditional on the sample, \( Y_t^*, Y_{t+1}^*, \ldots, Y_{T-1}^* \), are iid uniform, for \( t = 2, \ldots, T-1 \). Analogously, \( Y_{jt}^*, Y_{jt+1}^*, \ldots, Y_{jt+1}^* \), conditional on the (simulated) sample, is also iid uniform. Thus, given A1-A3, and by the basic properties of the block bootstrap (Künsch (1989), theorem 3.5), we also have that pointwise in \( u \):

\[ \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{ Y_t^* \leq u \} - F_j(u) \right) = \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{ Y_t \leq u \} - F_j(u) \right) + o_{P^*}(1), \quad \Pr - P_0, \quad (25) \]

and as \( l_T/T^{1/2} \to 0 \):

\[ \frac{1}{T} \sum_{t=1}^{T} \left( 1 \{ Y_t^* \leq u \} - F_j(u) \right) = \mu_{F_j}(u) + o_{P^*}(1), \quad \Pr - P_0. \quad (26) \]

Uniform convergence on \( U \) follows because of the stochastic equicontinuity of the left hand side of (25) and (26). Also, by the same argument used in Lemma A4 of Goncalves and White (2004: GW), and recalling A3(iv), it follows that for \( j = 1, \ldots, m \):

\[ P^*_j \left( \sup_{\theta_j \in \Theta_j} \sup_{u \in U} \left| \frac{1}{s} \sum_{n=1}^{s} f_j \left( u - (Y_{j,n}^*(\theta_j) - Y_{j,n}(\theta_j^*)) \right) \nabla_{\theta_j} Y_{j,n}^*(\theta_j) - \mu_{F_j}(u) \right| > \varepsilon \right) \to 0, \quad \Pr - P_j, \quad (27) \]

Thus, given (25)-(27), and using arguments similar to those used in the proof of Proposition 1, we see that after some elementary manipulations \( \sqrt{T} Z_{j,T,S}^*(u) \) can be written as:

\[ \sqrt{T} Z_{j,T,S}^*(u) = -\frac{2}{\sqrt{T}} \sum_{t=1}^{T} \left( 1 \{ Y_t^* \leq u \} - 1 \{ Y_t \leq u \} \right) (F_1(u) - F_j(u)) \]
By Theorem 2.2 in GW, \( \sqrt{T}(\hat{\theta}_{j,T}^* - \hat{\theta}_j) \) has the same limiting distribution as \( \sqrt{T}(\hat{\theta}_{j,T}^* - \theta_j^\dagger) \), c.z.p. Thus, by Theorem 3.5 in Künsch (1989), we see that pointwise in \( u \), \( Z_{j,T,S}(u) \) has the same limiting distribution as \( Z_{j,T,S}(u) \), c.z.p \( P_0 \) and \( P_j, j = 1, \ldots, m \) probability measure, respectively. By the Cramer Wold device, the same holds for any finite set of points in \( U \). Finally, by the empirical version of the block bootstrap (see e.g. Naik-Nimbalkar and Rajarshi (1994, Theorem 2.1)), it follows that \( \frac{1}{\sqrt{T}} \sum_{t=1}^T (1\{Y_t^* \leq u\} - 1\{Y_t \leq u\}) \) has the same limiting distribution as \( \frac{1}{\sqrt{T}} \sum_{t=1}^T (1\{Y_t \leq u\} - F_0(u)) \), as a process on \( U \), c.z.p \( P_0 \)-probability measure. Also, \( \frac{1}{\sqrt{S}} \sum_{n=1}^S (1\{Y_{j,n}^\dagger (\theta_j^\dagger) \leq u\} - 1\{Y_{j,n} (\theta_j^\dagger) \leq u\}) \) has the same limiting distribution as \( \frac{1}{\sqrt{S}} \sum_{n=1}^S (1\{Y_{j,n} \leq u\} - F_j(u)) \), as a process on \( U \), conditional on the simulated sample and for all simulated samples except a set of zero \( P_j \)-probability measure, \( j = 1, \ldots, m \).

The statement in part (i) of the theorem then follows as a straightforward consequence of the continuous mapping theorem.

(ii) The proof follows as a special case of part (i).
7 References


### Table 1: Comparison of Three Model Parameterizations Using Two Variables

<table>
<thead>
<tr>
<th>S, I</th>
<th>( \sqrt{T}Z_{T,S} )</th>
<th>Critval ((Z^*_{T,S}))</th>
<th>% &lt; 0</th>
<th>Min</th>
<th>Max</th>
<th>Stdevr</th>
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<tbody>
<tr>
<td></td>
<td>5%</td>
<td>10%</td>
<td>5%</td>
<td>10%</td>
<td>Sim1</td>
<td>Sim2</td>
</tr>
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<td>T,11</td>
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<td>0.844</td>
<td>0.109</td>
<td>0.726</td>
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<td>T,12</td>
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<td>0.065</td>
<td>0.0472</td>
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**Part I: Technology Shocks Assumed Normally Distributed; \( \alpha = 0.90, 0.95, 0.99; \) Benchmark sets \( \beta = 0.99 \)**

**Notes:** Results are based on examination of a version of the Christiano (1988) model. Simulated output and hours growth rates are compared. The first five columns of numerical entries report test statistic values (column 1) as well as 5% and 10% bootstrap critical values constructed using \( \sqrt{T}Z_{T,S} \) and \( \sqrt{T}Z^*_{T,S} \) (i.e. \( T/S > 0 \) and \( T/S \rightarrow 0 \), respectively, where \( T \) denotes the actual data sample size and \( S \) is the length of the simulated sample). For the case where \( S = T \), we set \( S = T \); say, we set \( S \) equal to \( \omega \) times the corresponding value of \( li \) when \( S = T \). All statistics are based on use of 20 uniformly distributed values for \( a \) across the range of the actual data; and bootstrap empirical distributions are constructed using 100 bootstrap replications. In addition to test statistics and critical values, various summary statistics constructed using the simulated data are given in the remaining columns of the table, for simulated output growth. Corresponding values based on the actual data (70 annual U.S. per capita output growth observations for the period 1933-2002) are: \% < 0 = 21.3; \( min = -0.128; \) \( max = 0.158; \) and \( stdevr = 0.044 \). We illustrate that the test has power against unreasonable alternatives in Part II. Under the benchmark scenario (Sim1), the errors driving the model are assumed to follow Student’s \( t \) distributions with unit variance, while under the alternatives, the errors are normally distributed (with variances 0.039 and 0.077 in Sim2 - see Christiano (1988) - and with variances estimated as discussed above in Sim3). See above for further details.
Table 2: Comparison of Three Model Parameterizations Using Four Variables

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<th>Sim2</th>
<th>Sim3</th>
<th>Sim1</th>
<th>Sim2</th>
<th>Sim3</th>
<th>Max Sim1</th>
<th>Sim2</th>
<th>Sim3</th>
<th>Stderr Sim1</th>
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Notes: See notes to Table 1. Simulated data being compared are output and hours growth rates, as well as the first lags of each of these variables. See above for further details.