

INFERENCE WHEN A NUISANCE PARAMETER IS NOT IDENTIFIED UNDER THE NULL HYPOTHESIS

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Many econometric testing problems involve nuisance parameters which are not identified under the null hypotheses. This paper studies the asymptotic distribution theory for such tests. The asymptotic distributions of standard test statistics are described as functionals of *chi-square processes*. In general, the distributions depend upon a large number of unknown parameters. We show that a transformation based upon a conditional probability measure yields an asymptotic distribution free of nuisance parameters, and we show that this transformation can be easily approximated via simulation. The theory is applied to threshold models, with special attention given to the so-called *self-exciting threshold autoregressive model*. Monte Carlo methods are used to assess the finite sample distributions. The tests are applied to U.S. GNP growth rates, and we find that Potter's (1995) threshold effect in this series can be possibly explained by sampling variation.

KEYWORDS: Asymptotic theory, nonlinear models, thresholds, identification, *p*-values, hypothesis testing.

1. INTRODUCTION

THIS PAPER STUDIES THE PROBLEM of inference in the presence of nuisance parameters which are not identified under the null hypothesis. The models considered take the form of additive nonlinearity, allowing for stochastic regressors and weak dependence. The asymptotic distributions of standard tests are nonstandard and nonsimilar, which means (among other things) that tabulation of critical values is impossible. This paper proposes a conditional transformation which is analogous to an asymptotic *p*-value, and has an asymptotic uniform distribution under the null hypothesis. The transformation is not directly calculable, but can be approximated using simple simulation techniques.

There are many econometric hypotheses of interest with unidentified nuisance parameters. Some examples include: (i) common (canceling) ARMA roots; (ii) no ARCH effect in a GARCH or ARCH-M model; (iii) constancy of a regression coefficient with the alternative that it follows an AR(1) process; (iv) a single regime or state against the alternative of multiple regimes. For brevity, this paper does not explicitly examine these cases, although the methods of this paper can be extended to handle these cases as well.²

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²An earlier version of this paper (Hansen (1991)) used a more general notational framework which included nonlinear maximum likelihood estimation.

This paper concentrates on regression models with additive nonlinearity, which take the form $y_t = x_t' \alpha + h(z_t, \gamma)' \theta + \varepsilon_t$. Here, the null hypothesis of interest is whether the nonlinear term $h(z_t, \gamma)$ enters the regression. When $\theta = 0$ the parameter γ is not identified. Specific examples of additive nonlinearity include the following. (i) Box-Cox transformations: $h(z_t, \gamma) = (z_t^\gamma - 1)/\gamma$; (ii) structural change: $h(z_t, \gamma) = \{t/n \leq \gamma\} z_t$, where $\{\cdot\}$ is the indicator function; (iii) threshold models: $h(z_t, \gamma) = \{z_t \leq \gamma\} z_t$; (iv) Bierens's (1990) consistent tests of functional form: $h(z_t, \gamma) = \exp(\gamma' z_t)$; (v) White's (1989) neural network tests of functional form: $h(z_t, \gamma) = \psi(\gamma' z_t)$, where $\psi(\cdot)$ is the logistic function.

There have been several recent papers on the subject of unidentified nuisance parameters which examine related issues. Andrews (1993b) analyzed tests for structural change. Andrews and Ploberger (1994) explore optimal testing but do not discuss methods to obtain critical values in practice. Andrews (1993a) discusses a range of econometric examples which suffer from the problem of unidentified nuisance parameters. King and Shively (1991) discuss the merits and difficulties of reparameterization as a means to handle the problem. Stinchcombe and White (1993) examine White's neural network tests. All of these papers investigate different aspects of the problem and should be viewed as complementary to the results of this paper.

In the next section, we introduce additive nonlinear regression and discuss pointwise and global test statistics. Section 3 introduces the conditional p -value transformation. Section 4 examines threshold regression, including the self-exciting threshold autoregressive (SETAR) model of Tong (1983). A Monte Carlo study is reported in Section 5. We find that the LM-based tests have excellent size and good power. In Section 6, we apply these tests to Potter's (1995) SETAR model of U.S. GNP and find that the apparent "significant" threshold effect may be explained by sampling error. Section 7 contains a brief conclusion. Proofs are left to an Appendix. Concerning notation, let $|A| = (\text{tr}(A'A))^{1/2}$ denote the Euclidean norm of a matrix A , let $\|A\|_r = (E|A|^r)^{1/r}$ denote the L^r -norm of a random matrix, and let \Rightarrow denote weak convergence with respect to the uniform metric.

2. FRAMEWORK AND TEST STATISTICS

The data is $\{w_t = (y_t, x_t) : t = 1, \dots, n\}$, which is a draw from some underlying probability space, and satisfies the regression relationship $y_t = x_{1t}' \beta_1 + h(x_t, \gamma)' \beta_2 + \varepsilon_t$, where x_{1t} is a $k_1 \times 1$ subvector of the k vector x_t , and ε_t is a real-valued martingale difference sequence with respect to some increasing set of sigma-fields \mathcal{F}_t to which ε_t and x_{t+1} are adapted, and $E\varepsilon_t^2 = \sigma^2 < \infty$. γ takes values in Γ , a bounded subset of R^q , and the function $h(\cdot, \cdot)$ maps $R^k \times \Gamma$ into R^m . We will typically write $h(x_t, \gamma)$ more simply as $h_t(\gamma)$. It will be convenient to write the model in the alternate form

$$(1) \quad y_t = x_t(\gamma)' \beta + \varepsilon_t$$

where $x_t(\gamma) = (x_{1t}', h_t(\gamma)')'$ and $\beta = (\beta_1' \beta_2)'$.

The question of interest is whether the nonlinear term $h_t(\gamma)$ enters the regression, that is, whether $\beta_2 = 0$. The distributional theory will be facilitated by a local-to-null reparameterization: $\beta_2 = c/\sqrt{n}$. The null hypothesis is $H_0 : c = 0$ with alternative $H_1 : c \neq 0$. The test is nonstandard since γ does not enter the regression (and therefore is not identified) under H_0 .

Under H_0 the model simplifies to $y_t = x'_{1t} \beta_1 + \varepsilon_t$. The OLS estimators are $\tilde{\beta}_1 = (\sum_{t=1}^n x_{1t} x'_{1t})^{-1} (\sum_{t=1}^n x_{1t} y_t)$ and $\tilde{\sigma}_n^2 = \sum_{t=1}^n \tilde{\varepsilon}_t^2 / (n - k_1)$, where $\tilde{\varepsilon}_t = y_t - x'_{1t} \tilde{\beta}_1$.

Under H_1 , if γ were known, then β could be estimated by ordinary least squares on (1), yielding estimates $\hat{\beta}(\gamma) = (\sum_{t=1}^n x_t(\gamma) x_t(\gamma)')^{-1} (\sum_{t=1}^n x_t(\gamma) y_t)$, residuals $\hat{\varepsilon}_t(\gamma) = y_t - x_t(\gamma)' \hat{\beta}(\gamma)$, and sample variance $\hat{\sigma}_n^2(\gamma) = \sum_{t=1}^n \hat{\varepsilon}_t(\gamma)^2 / (n - (k_1 + m))$. When γ is unknown, then its least-squares estimate can be found by minimization of $\hat{\sigma}_n^2(\gamma)$ over $\gamma \in \Gamma$, yielding $\hat{\gamma} = \text{argmin } \hat{\sigma}_n^2(\gamma)$ and $\hat{\beta} = \hat{\beta}(\hat{\gamma})$. These are also the MLE when ε_t is iid Gaussian. It will be useful to define the regression scores $s_t(\gamma) = x_t(\gamma) \varepsilon_t$ and their estimates under H_0 and H_1 , respectively, $\hat{s}_t(\gamma) = x_t(\gamma) \hat{\varepsilon}_t(\gamma)$ and $\tilde{s}_t(\gamma) = x_t(\gamma) \tilde{\varepsilon}_t$.

If γ were known, then the testing problem would not be complicated. A heteroskedasticity-robust Wald test takes the form

$$T_n(\gamma) = n \hat{\beta}(\gamma)' R (R \hat{V}_n^*(\gamma) R)^{-1} R' \hat{\beta}(\gamma)$$

where R is the selector matrix $R = (0 I_p)'$, $\hat{V}_n^*(\gamma) = M_n(\gamma, \gamma)^{-1} \hat{V}_n(\gamma) M_n(\gamma, \gamma)^{-1}$, $\hat{V}_n(\gamma) = \frac{1}{n} \sum_{t=1}^n \hat{s}_t(\gamma) \hat{s}_t(\gamma)'$, and $M_n(\gamma_1, \gamma_2) = \frac{1}{n} \sum_{t=1}^n x_t(\gamma_1) x_t(\gamma_2)'$. Alternatively, a Lagrange multiplier statistic is found by setting $\tilde{V}_n(\gamma) = \frac{1}{n} \sum_{t=1}^n \tilde{s}_t(\gamma) \tilde{s}_t(\gamma)'$. Tests which assume homoskedastic errors can be formed similarly.

If γ were known *a priori*, then under conventional regularity conditions $T_n(\gamma)$ would have a “point-optimal” interpretation and an approximate χ_m^2 null distribution in large samples. Hence H_0 can be tested by selecting a “reasonable” value of γ a priori. We might expect this test procedure to work well if γ is *known* a priori, or if the selected test statistic is not sensitive to the choice of γ . For many of the examples of interest, however, neither condition holds. The researcher is left with an unpleasant dilemma. Either γ is selected in a completely arbitrary way (and thereby sacrifices power) or γ is selected in some data-dependent fashion, in which case the chi-square distributional approximation will be invalid. In addition, researchers who estimate unrestricted models will be generating estimates of γ . They will be interested in the question: “Does the unrestricted model fit statistically better than the restricted model?” which cannot be answered by a point-optimal test. To avoid these difficulties, we turn to tests which do not require prior knowledge of γ .

Davies (1977, 1987) suggested testing H_0 by $\sup T_n = \sup_{\gamma \in \Gamma} T_n(\gamma)$, which equals the LR statistic when $T_n(\gamma)$ is the pointwise likelihood ratio statistic. Andrews and Ploberger (1994) examine tests of H_0 and argue that superior local power can be constructed from the statistics $\text{ave} T_n = \int_{\Gamma} T_n(\gamma) dW(\gamma)$ and $\exp T_n = \ln(\int_{\Gamma} \exp(\frac{1}{2} Z(\gamma)) dW(\gamma))$. All three statistics can be written as functions $g(T_n)$, where $g(\cdot)$ maps functionals on Γ to R , and we write $T_n = \{T_n(\gamma) : \gamma \in \Gamma\}$ as a random function on Γ . Each function g is continuous with respect to the

uniform metric, monotonic in the sense that if $Z_1(\gamma) \leq Z_2(\gamma)$ for all γ then $g(Z_1) \leq g(Z_2)$, and has the property that if $Z(\gamma) \rightarrow \infty$ for γ for some subset of Γ with positive W -measure, then $g(Z) \rightarrow \infty$. We will denote the test statistic as $g_n = g(T_n)$ for any choice of g .

Define $M(\gamma_1, \gamma_2) = E(x_t(\gamma_1)x_t(\gamma_2)')$ and $\bar{x}_t = \sup_{\gamma \in \Gamma} |x_t(\gamma)|$.

ASSUMPTION 1: w_t is strictly stationary and absolutely regular with mixing coefficients $\eta(m) = O(m^{-A})$ for some $A > v/(v - 1)$ and $r \geq v > 1$; $E|\bar{x}_t|^{4r} < \infty$; $E|\varepsilon_t|^{4r} < \infty$; and $\inf_{\gamma \in \Gamma} \det(M(\gamma, \gamma)) > 0$.

The absolute regular mixing coefficient $\eta(\mathcal{A}, \mathcal{B})$ between σ -fields \mathcal{A} and \mathcal{B} is defined as

$$\eta(\mathcal{A}, \mathcal{B}) = \frac{1}{2} \sup_{(i,j) \in (I,J)} \sum |P(A_i \cap B_j) - P(A_i)P(B_j)|,$$

where $A_i \subset \mathcal{A}$, $B_j \subset \mathcal{B}$, and the supremum is taken over all the finite partitions $(A_i)_{i \in I}$ and $(B_j)_{j \in J}$ respectively \mathcal{A} and \mathcal{B} measurable. Absolute regularity was first defined by Volkonskii and Rozanov (1959), and is stronger than strong mixing yet weaker than uniform mixing. Pham and Tran (1985) have shown that a wide class of linear processes with iid innovations (such as ARMA processes) are absolutely regular when the innovation has a bounded, continuous density, and thus these processes satisfy Assumption 1. Define $K_n(\gamma_1, \gamma_2) = \frac{1}{n} \sum_{t=1}^n s_t(\gamma_1)s_t(\gamma_2)'$, and $K(\gamma_1, \gamma_2) = E(s_t(\gamma_1)s_t(\gamma_2)')$. Under the stated moment conditions, we see that for all $\gamma_1, \gamma_2 \in \Gamma$, $M_n(\gamma_1, \gamma_2) \rightarrow M(\gamma_1, \gamma_2)$ and $K_n(\gamma_1, \gamma_2) \rightarrow K(\gamma_1, \gamma_2)$ a.s. We will need the stronger requirement of uniform convergence. At this point we give high-level conditions, supplying primitive conditions in Section 4.

ASSUMPTION 2: For some $B < \infty$ and $\lambda > 0$, $\|(h_t(\gamma) - h_t(\gamma'))\varepsilon_t\|_{2v} \leq B|\gamma - \gamma'|^\lambda$.

ASSUMPTION 3: $M_n(\gamma_1, \gamma_2)$ and $K_n(\gamma_1, \gamma_2)$ converge almost surely to $M(\gamma_1, \gamma_2)$ and $K(\gamma_1, \gamma_2)$, respectively, uniformly over $\gamma_1, \gamma_2 \in \Gamma$.

Let $S_n(\gamma) = (1/\sqrt{n})\sum_{t=1}^n s_t(\gamma)$, and let $S(\gamma)$ denote a mean zero Gaussian process with covariance kernel $K(\gamma_1, \gamma_2)$. This means that for any $\{\gamma_1, \dots, \gamma_k\} \in \Gamma^k$, $\{S(\gamma_1), \dots, S(\gamma_k)\}$ is multivariate normal with mean zero and covariances $E(S(\gamma_j)S(\gamma_l)') = K(\gamma_j, \gamma_l)$. Next, let $\bar{S}(\gamma) = R'M(\gamma, \gamma)^{-1}S(\gamma)$ which is a mean-zero Gaussian process with covariance kernel

$$\bar{K}(\gamma_1, \gamma_2) = R'M(\gamma_1, \gamma_1)^{-1}K(\gamma_1, \gamma_2)M(\gamma_2, \gamma_2)^{-1}R.$$

THEOREM 1: Under Assumptions 1-3, $S_n \Rightarrow S$, $T_n \Rightarrow T^c$, and $g_n \Rightarrow g^c = g(T^c)$, where

$$T^c(\gamma) = (\bar{S}(\gamma)' + c' \bar{Q}(\gamma)') \bar{K}(\gamma, \gamma)^{-1} (\bar{S}(\gamma) + \bar{Q}(\gamma)c)$$

and $\bar{Q}(\gamma) = R' M(\gamma, \gamma)^{-1} M(\gamma, \gamma_0) R$, with γ_0 the true value of γ when $c \neq 0$.

Theorem 1 gives the asymptotic distribution of the test statistic sequence T_n and the global test g_n under the local alternative $\beta_2 = c/\sqrt{n}$. Thus the asymptotic null distribution of T_n is $T^0(\gamma) = \bar{S}(\gamma)' \bar{K}(\gamma, \gamma)^{-1} \bar{S}(\gamma)$, which has for each $\gamma \in \Gamma$ a marginal chi-square distribution. Hence, we call T^0 a *chi-square process*, and similarly T^c a *noncentral chi-square process*. Since the null distribution of $g^0 = g(T^0)$ depends, in general, upon the covariance function \bar{K} , critical values cannot be tabulated except in special cases.³

3. THE P-VALUE TRANSFORMATION

Let $F^0(\cdot)$ denote the distribution function of g^0 , and define $p_n = 1 - F^0(g_n)$. Tests based on g_n and p_n are equivalent since F^0 is monotonic and continuous. From Theorem 1 we see that $p_n \Rightarrow p^c$, where $p^c = 1 - F^0(g^c)$. In particular, the null distribution is $p^0 \sim U$, the uniform distribution on $[0, 1]$. Thus the asymptotic null distribution of p_n is free of nuisance parameters. Our test is to reject H_0 if $p_n \leq \alpha$. When the exact null distribution of p_n is U , then α represents the size of the test; otherwise α is an asymptotic size, since

$$(2) \quad \lim_{n \rightarrow \infty} P\{p_n \leq \alpha | c = 0\} = \alpha.$$

We call p_n an “asymptotic p -value.”

Strictly speaking, the label “ p -value” and “asymptotic size” are correct only if the convergence in (2) is uniform⁴ over the null hypothesis. Note, however, that the distributions of T_n , g_n , and p_n are invariant to the parameters β_1 and σ^2 . Thus in the special case in which x_t is strictly exogenous and ε_t is i.i.d. normal, the distribution of p_n is free of nuisance parameters, so H_0 is simple and the convergence in (2) is trivially uniform. In the general regression context, uniform convergence may not hold. This is common in econometric testing problems and will not be emphasized in this paper.

The asymptotic power function associated with the test is

$$\pi_\alpha(c) = \lim_{n \rightarrow \infty} P\{p_n \leq \alpha | c\} = P\{F^0(g^c) \geq 1 - \alpha | c\}.$$

It is possible to show that as $|c|$ becomes large, $g^c \rightarrow \infty$ almost surely, so $p^c \rightarrow 0$ a.s. Hence for any $\alpha > 0$, $\pi_\alpha(c) \rightarrow 1$ as $|c| \rightarrow \infty$, and the test has nontrivial local power.

³For example, when testing for structural change of unknown timing (Andrews (1993b)) the asymptotic distribution only depends upon Γ and m .

⁴For a definition see Sweeting (1980).

The random variable g^0 can be written as a continuous functional of the Gaussian process $S(\gamma)$, which is completely described by its covariance kernel $K(\gamma_1, \gamma_2)$. We can construct two estimates of this kernel, derived by the Wald and LM principles: $\hat{K}_n(\gamma_1, \gamma_2) = \frac{1}{n} \sum_{t=1}^n \hat{s}_t(\gamma_1) \hat{s}_t(\gamma_2)'$ and $\tilde{K}_n(\gamma_1, \gamma_2) = \frac{1}{n} \sum_{t=1}^n \tilde{s}_t(\gamma_1) \tilde{s}_t(\gamma_2)'$. One attraction of the estimator \hat{K}_n is that, like an LM statistic, it only involves estimation under the null hypothesis.

Now operate conditionally on the sample w . Denote by \hat{S}_n a conditional mean-zero Gaussian process with covariance kernel $\hat{K}_n(\cdot, \cdot)$, by \hat{T}_n a conditional chi-square process with covariance kernel $R' M_n(\gamma_1, \gamma_1)^{-1} \hat{K}_n(\gamma_1, \gamma_2) M_n(\gamma_2, \gamma_2)^{-1} R$, and let $\hat{g}_n = g(\hat{T}_n)$. Let \hat{F}_n denote the conditional distribution function of \hat{g}_n , conditional on the sample w , and set $\hat{p}_n = 1 - \hat{F}_n(g_n)$. Similarly, let \tilde{S}_n denote a conditional mean-zero Gaussian process with covariance kernel $\tilde{K}_n(\cdot, \cdot)$, and similarly $\tilde{T}_n, \tilde{g}_n, \tilde{F}_n$, and \tilde{p}_n .

These are not vacuous definitions. \hat{S}_n and \hat{T}_n can be generated by letting $\{v_t\}_{t=1}^n$ be i.i.d. $N(0, 1)$ random variables, and setting

$$(3) \quad \hat{S}_n(\gamma) = \frac{1}{\sqrt{n}} \sum_{t=1}^n \hat{s}_t(\gamma) v_t$$

and

$$(4) \quad \hat{T}_n(\gamma) = \hat{S}_n(\gamma)' M_n(\gamma, \gamma)^{-1} R (R' \hat{V}_n^*(\gamma) R)^{-1} R' M_n(\gamma, \gamma)^{-1} \hat{S}_n(\gamma).$$

Similarly, \tilde{S}_n and \tilde{T}_n can be generated by

$$(5) \quad \tilde{S}_n(\gamma) = \frac{1}{\sqrt{n}} \sum_{t=1}^n \tilde{s}_t(\gamma) v_t$$

and

$$(6) \quad \tilde{T}_n(\gamma) = \tilde{S}_n(\gamma)' M_n(\gamma, \gamma)^{-1} R (R' \hat{V}_n^*(\gamma) R)^{-1} R' M_n(\gamma, \gamma)^{-1} \tilde{S}_n(\gamma).$$

THEOREM 2: *Under Assumptions 1 and 2, $\hat{p}_n = p_n + o_p(1)$ and $\tilde{p}_n = p_n + o_p(1)$. Hence $\hat{p}_n \Rightarrow p^c$ and $\tilde{p}_n \Rightarrow p^c$, and under H_0 the asymptotic distribution of both \hat{p}_n and \tilde{p}_n is U .*

Theorem 2 shows that \hat{p}_n and \tilde{p}_n are asymptotically equivalent to p_n , under both the null hypothesis and local alternative $\beta_2 = c/\sqrt{n}$. The proof of the Theorem runs roughly as follows. We show that $\hat{S}_n \Rightarrow_p S$, where “ \Rightarrow_p ” denotes “weak convergence in probability” as defined by Gine and Zinn (1990). This implies that $\hat{g}_n \Rightarrow_p g^0$ and hence $\hat{F}_n(x) \rightarrow_p F^0(x)$, uniformly in x . Thus $\hat{p}_n = 1 - \hat{F}_n(g_n) + o_p(1)$, as stated. The proof has similarities to those used to derive asymptotic properties of the bootstrap.

The conditional distribution functions $\hat{F}_n(\cdot)$ and $\tilde{F}_n(\cdot)$ are not directly observed, so neither are the random variables \hat{p}_n and \tilde{p}_n . We can approximate either \hat{F}_n or \tilde{F}_n to any desired degree of accuracy, however, using standard

simulation techniques. Equations (3)–(6) show that a random draw from either conditional distribution can be made using a sample of iid $N(0, 1)$ variables. Take, for example, \hat{F}_n . For $j = 1, \dots, J$, execute the following steps:

(i) generate $\{v_{ij}\}_{i=1}^n$ iid $N(0, 1)$ random variables (using a random number generator);

(ii) set $S_n^j(\gamma) = (1/\sqrt{n})\sum_{i=1}^n \hat{s}_i(\gamma)v_{ij}$;

(iii) set $T_n^j(\gamma) = S_n^j(\gamma)'M_n(\gamma, \gamma)^{-1}R(R'\hat{V}_n^*(\gamma)R)^{-1}R'M_n(\gamma, \gamma)^{-1}S_n^j(\gamma)$;

(iv) set $g_n^j = g(T_n^j)$.

This gives a random sample (g_n^1, \dots, g_n^J) of J observations from the conditional distribution \hat{F}_n . Then compute the percentage of these artificial observations which exceed the actual test statistic $g_n: \hat{p}_n^J = \frac{1}{J}\sum_{j=1}^J \{g_n^j \geq g_n\}$. By the Glivenko-Cantelli Theorem, for any sample $\hat{p}_n^J \rightarrow_p 1 - \hat{F}_n(g_n) = \hat{p}_n$ as $J \rightarrow \infty$. Since J is under the control of the econometrician, \hat{p}_n^J can be made arbitrarily close to \hat{p}_n by picking sufficiently large J , and thus can be used as our test statistic in place of \hat{p}_n . Analogously, we can construct \tilde{p}_n^J , using $S_n^j(\gamma) = (1/\sqrt{n})\sum_{i=1}^n x_i(\gamma)\tilde{\varepsilon}_i v_{ij}$ instead in step (ii). To select J in practice, an appeal to the central limit theorem gives an asymptotic standard error for \hat{p}_n^J of $\sqrt{\hat{p}_n^J(1 - \hat{p}_n^J)/J}$. For example, when $\hat{p}_n^J = 0.05$, setting $J = 1000$ yields a standard error of only 0.007.

When the regression error is conditionally homoskedastic: $E(\varepsilon_t^2 | \mathcal{F}_{t-1}) = \sigma^2$ a.s., the asymptotic expressions are significantly simpler. To approximate the asymptotic p -value using this information, for $j = 1, \dots, J$, set $S_n^j(\gamma) = (1/\sqrt{n})\sum_{i=1}^n x_i(\gamma)v_{ij}$, where v_{ij} are iid $N(0, 1)$ draws from a random number generator,

$$T_n^j(\gamma) = S_n^j(\gamma)'M_n(\gamma, \gamma)^{-1}R(R'M_n(\gamma, \gamma)^{-1}R)^{-1}R'M_n(\gamma, \gamma)^{-1}S_n^j(\gamma),$$

and $\hat{p}_n^J = (1/J)\sum_{j=1}^J \{g(T_n^j) \geq g_n\}$. The same arguments as before show that $\hat{p}_n^J \rightarrow_p \hat{p}_n$ as $J \rightarrow \infty$ and $\hat{p}_n \rightarrow_d p^c$ as $n \rightarrow \infty$, with the null distribution $p^0 \sim U$.

If Γ is a continuous parameter space, calculation of g_n and replications g_n^j might be excessively costly. In this case, it may be reasonable to replace Γ by a discrete approximation $\Gamma_A = (\gamma_1, \dots, \gamma_A)$. Then the actual test statistic is $g_{An} = g(T_n, \Gamma_A)$, where the notation is explicit about the dependence upon the region Γ_A . The simulated draws are $\hat{g}_{An}^j = g(\hat{T}_n^j, \Gamma_A)$ and the p -values are $\hat{p}_{An}^J = (1/J)\sum_{j=1}^J \{\hat{g}_{An}^j \geq g_{An}\}$. The null asymptotic theory goes through unaffected. Specifically, $g_{An} \Rightarrow g(T, \Gamma_A) = g_A$, say, and $\hat{p}_{An}^J \rightarrow_d p_A^c$, with $p_A^0 \sim U$. The power of the test may be adversely affected, however, if the selected approximation Γ_A is insufficiently dense in Γ . Essentially, the function $T_n(\gamma)$ may have a sharp peak at some $\gamma_n \in \Gamma$ which is missed by the set Γ_A . A general solution to this problem may be impossible, since the appropriate choice of Γ_A will depend upon the smoothness of $T_n(\gamma)$. If this is the case, several choices of Γ_A could be used to assess sensitivity. In some applications, such as the threshold models explored in the next section, this is not an issue, as the sample function $T_n(\gamma)$ are naturally step functions with known step-points.

4. THRESHOLD REGRESSION

A typical threshold regression model takes the form (1) with $h(x_t, \gamma) = \{z_t \leq \gamma\}x_t$, where $\{\cdot\}$ denotes the indicator function, z_t is an element of x_t , $x_{1t} = x_t$, and $\Gamma = [\gamma_L, \gamma_U]$. Threshold models are quite common in applied econometrics, typically interpreted as sample splits. By formally treating γ as an unknown parameter, the threshold model allows the selection of γ to be made conditional on the data.

The regression function and all test statistics are functions of z_t and γ only through the indicator function $\{z_t \leq \gamma\}$. For any monotonic transformation $Z_n(\cdot): \Gamma \rightarrow \mathcal{T}$, we have $\{z_t \leq \gamma\} = \{z_{nt} \leq \tau\}$ where $z_{nt} = Z_n(z_t)$ and $\tau = Z_n(\gamma)$. An ideal choice for $Z_n(\cdot)$ is the empirical distribution function of z_t . One advantage is that the transformed threshold parameter τ is unit-free and lies in the generic region $[0, 1]$. The pointwise test statistics are ill-behaved for extreme values of τ , so \mathcal{T} should not include values of τ too close to 0 or 1. There is no obvious criteria by which to make this choice, but a similar dilemma appears in the changepoint literature, where recently Andrews (1993b) suggested $\mathcal{T} = [.15, .85]$.

Davies (1977) mentioned threshold models as a possible application of his distributional theory, but did not investigate whether his conditions were satisfied. In fact, they are not. Davies assumed that the limit process $T(\gamma)$ has a derivative, but this is not the case in threshold models. Take the special case where $x_t = 1$ and $z_t = y_{t-1}$. Here, the asymptotic process $T(\gamma)$ is known to be a normalized squared Brownian bridge (see Chan (1990)), which is nowhere differentiable and thus has infinite total variation. The bound Davies uses is a function of the sample total variation, which is finite in any sample yet diverges to infinity as $n \rightarrow \infty$, rendering an asymptotic size of zero. Table I compares Davies' procedure with the test based on the asymptotic critical values, where the innovations ε_t are iid $N(0, 1)$, and the sample size is varied from 50 to 1000. The null sets $\theta = 0$ and the alternative sets $\theta = .5$ and $\gamma = 0$. The first five columns show the rejection frequency under the null, in which case Davies' suggested procedure is extremely conservative. The cost shows up in the rejection frequency under the alternative (the final five columns), where the Davies' procedure yields tests with power less than the nominal 5% rate for samples of 500 or less. As expected from the asymptotic theory, Davies' bound is not useful in this context.

TABLE I
SIZE AND POWER OF NOMINAL 5% SIZE TESTS FOR SHIFT IN MEAN

Sample Size	Null $\theta = 0$					Alternative $\theta = .5$				
	50	100	250	500	1000	50	100	250	500	1000
Davies Procedure	.00	.01	.01	.01	.00	.01	.02	.03	.05	.09
Asymptotic Criticals	.02	.04	.04	.04	.05	.04	.08	.12	.19	.34

Instead, we can apply the conditional transformation of Section 3 to obtain appropriate asymptotic critical values. The following theorem shows that this procedure is justified.

THEOREM 3: *Assume that Assumption 1 holds with $r > v$, and z_t has density function $f(z)$ such that $\sup_{x \in R} f(x) = \bar{f} < \infty$. Then Assumptions 2 and 3 are satisfied, and hence Theorems 1 and 2 hold.*

The threshold model assumes that the threshold variable z_t is known a priori. This is not always the case. In many examples, all that is known is that z_t is some element of x_{2t} , some $k_2 \times 1$ subvector of x_t . In this case, we can write $z_t = x_{2t}(d)$, the d th element of x_{2t} . In this notation, we can think of d as a parameter, taking values in the discrete set $D = \{1, 2, \dots, k_2\}$. The standard threshold model emerges as the special case in which d is known a priori. In the threshold selection model d is unknown and must be estimated along with the other parameters.

For fixed (γ, d) , the model is linear in (α, θ) so can be estimated by ordinary least squares. The global estimates of the parameters can be found by minimization of the resulting least squares variance estimate over $(\gamma, d) \in (\Gamma \times D)$. Under the null hypothesis both the parameters γ and d are not identified, so the pair (γ, d) is treated as we had treated γ in the earlier sections. Thus, the pointwise test statistics $T_n(\gamma, d)$ may be found for each (γ, d) , and the transformations g_n , \hat{p}_n , and \hat{p}'_n found as before, replacing the argument γ by (γ, d) .

It is not hard to see that the asymptotic theory is essentially unaffected. Since D is a finite set, all convergence results are uniform over $d \in D$.

COROLLARY 1: *If $z_t = x_{2t}(d)$, assume that Assumption 1 holds with $r > v$, and x_{2t} has density function $f(x)$ such that $\sup_{x \in R^{k_2}} f(x) = \bar{f} < \infty$. Then Assumptions 2 and 3 are satisfied, and hence Theorems 1 and 2 hold for the threshold selection model.*

A special example of a threshold regression is the so-called *self-exciting threshold autoregressive model* (SETAR), which has received considerable recent attention in the nonlinear time series literature. The model takes the form

$$(7) \quad y_t = \alpha_0 + \alpha_1 y_{t-1} + \dots + \alpha_p y_{t-p} + (\theta_0 + \theta_1 y_{t-1} + \dots + \theta_p y_{t-p}) \{y_{t-d} \leq \gamma\} + \varepsilon_t$$

with $\{\varepsilon_t, \mathcal{F}_t\}$ a MDS. The model (7) falls in the class of threshold selection models by setting $x_t = (1 \ y_{t-1} \ \dots \ y_{t-p})$ and $z_t(d) = y_{t-d}$, and is largely due to Tong (1983). Chan (1990) proved an analog of Theorem 1 under the assumption that d is known and ε_t is iid Gaussian. Chan (1991) approximated the tail of the asymptotic distribution for general p (but still fixed and known d). General results, allowing arbitrary distribution functions for ε_t , general functionals on Γ such as the Andrews-Ploberger optimal transformations, and treating d as an

unknown parameter, have not been treated before. Our theory easily handles these generalizations.

COROLLARY 2: Suppose that $\{y_t\}$ is generated by (7) with ε_t independent and identically distributed, all the roots of the characteristic equation $z^p - \alpha_1 z^{p-1} - \dots - \alpha_p = 0$ lie strictly within the unit circle, $E|\varepsilon_t|^{4r} < \infty$ for some $r > 1$, and the density of ε_t is bounded and continuous. Then under $H_0: \theta_0 = \dots = \theta_p = 0$, $\hat{p}_n \Rightarrow U$ and $\tilde{p}_n \Rightarrow U$.

The assumptions of Corollary 2 imply that y_t is strictly stationary and absolutely regular with exponentially declining mixing coefficients (see Pham and Tran (1985)), establishing Assumption 1. It also implies that y_t has a bounded density, completing the requirements for Corollary 1. The only reason why we technically restrict attention to the null hypothesis is because it has not been investigated whether or not SETAR processes are absolutely regular. Similarly, the restriction that ε_t is iid is made only to guarantee absolute regularity. It is quite likely that this condition is not necessary and could be replaced by a martingale difference condition. Hence while the assumptions include homoskedasticity, we do not think that it is prudent to impose this condition when constructing test statistics.

5. FINITE SAMPLE DISTRIBUTION

To assess the usefulness of the testing methodology, finite sample distributional results are reported in a simple Monte Carlo simulation study. We use the threshold-selection SETAR model (7) with $p = 1, 2$, and 3 , $\alpha_1 = .3$, $\alpha_0 = \alpha_2 = \alpha_3 = 0$, ε_t iid $N(0, 1)$ and two sample sizes, $n = 100$ and $n = 200$. When $p > 1$, both the threshold lag d and the threshold γ were estimated by least squares. The experiments were done for tests of size 10%, 5%, and 1%, but only those for size 5% are reported since no differences were observed.

The test statistics functions $T_n(\gamma, d)$ were calculated using four different covariance matrices: (i) standard Wald (W); (ii) standard Lagrange multiplier (LM); (iii) MacKinnon-White (1985) "jackknife" heteroskedasticity-consistent Wals (W^h); (iv) Eicker-White heteroskedasticity-consistent Lagrange multiplier (LM^h). From each of these test statistic functions, two functionals were used: the supremum (sup), and average (ave) (see Section 2) to generate the test statistics g_n and the simulated p -values \hat{p}_n^J . Due to the large computational requirements of the simulation design, the number of internal simulation replications was set at $J = 500$ and the region T for the calculation of the test statistics (see Section 4) was set at $T = [.2, .8]$. A ninth test statistic, included for comparison, is the S_3 test of Luukkonen, Saikkonen, and Terasvirta (1988), which is an LM-type test for a smooth transition autoregressive (STAR) model. The simulation study reported in Luukkonen, et al., showed that the S_3 test has good power.

TABLE II
FINITE SAMPLE SIZE OF ASYMPTOTIC 5% TESTS

AR order	$n = 100$			$n = 200$		
	$p = 1$	$p = 2$	$p = 3$	$p = 1$	$p = 2$	$p = 3$
SupW	.07	.07	.07	.05	.05	.08
AveW	.05	.05	.04	.04	.04	.06
SupLM	.05	.04	.03	.04	.04	.05
AveLM	.05	.04	.03	.04	.04	.06
SupW ^h	.14	.21	.32	.09	.13	.19
AveW ^h	.07	.09	.10	.05	.06	.07
SupLM ^h	.04	.02	.03	.05	.04	.02
AveLM ^h	.04	.04	.03	.04	.04	.05
STAR	.05	.04	.03	.04	.04	.04

We report in Table II the actual size (setting $\theta_0 = \theta_1 = 0$). For samples of size 100, reported in the first three columns, 2000 simulated samples were drawn. For samples of size 200, reported in the final three columns, 1000 simulated samples were used for $p = 1$ and $p = 2$, and 500 for $p = 3$. We find that the asymptotic approximation is excellent for the “standard” tests (without the heteroskedasticity correction). The heteroskedasticity-consistent LM tests are slightly conservative for large p , and the heteroskedasticity-consistent Wald statistic supW^h is excessively liberal, especially at $n = 100$. Based on these results, our recommendation is to avoid the supW^h statistic unless the sample size is very large.

To assess power, we consider two specifications, using first an intercept shift ($\theta_0 \neq 0$) and second a slope shift ($\theta_1 \neq 0$). In both specifications, γ is selected so that it is approximately the median of y_{t-1} . (Since there is no closed-form expression for the distribution or density of y_t , the median was calculated by simulation.) For each simulation design with $p \leq 2$, 1000 replications were made, and 500 replications were made when $p = 3$. Finite sample 5% critical values for the test statistics were calculated from the .05 percentile of the empirical distribution of the tests calculated under the null. For each p and each n , the power of the Wald and LM tests were nearly identical, so we only report in Table III the results of the LM-based tests.

As expected, power is increasing in $|\theta_0|$ or $|\theta_1|$, increasing in n , and decreasing with p . Indeed, the effect of the estimated model order (p) is quite strong, indicating that the cost of over-fitting is high. The theory of Andrews-Ploberger (1994) suggests that aveLM should be optimal against local alternatives, yet this is not supported by the simulation evidence. Instead, we find that for shifts in the intercept supLM has the best power, yet for shifts in the slope aveLM dominates. It is also interesting to compare the power of our tests with the STAR test of Luukkonen, et al. The STAR test does remarkably well against the shift in slope, although less well against the shift in the intercept. It also does better against local alternatives than distant alternatives, which is expected since it was derived using a local power argument. It is also possible to see a slight

TABLE III
FINITE SAMPLE POWER OF 5% SIZE TESTS

		<i>n</i> = 100				<i>n</i> = 200			
		θ_0		θ_1		θ_0		θ_1	
		-.75	-1.25	-.6	-1.0	-.75	-1.25	-.6	-1.0
<i>p</i> = 1	SupLM	.29	.70	.25	.69	.63	.99	.55	.98
	SupLM ^h	.28	.65	.22	.62	.61	.98	.52	.96
	AveLM	.21	.50	.38	.84	.39	.86	.67	.99
	AveLM ^h	.20	.47	.36	.80	.35	.83	.63	.98
	STAR	.18	.42	.31	.73	.34	.72	.58	.97
<i>p</i> = 2	SupLM	.22	.61	.18	.60	.29	.70	.25	.69
	SupLM ^h	.20	.55	.16	.52	.28	.65	.22	.62
	AveLM	.13	.27	.25	.67	.21	.50	.38	.84
	AveLM ^h	.12	.26	.21	.61	.20	.47	.35	.80
	STAR	.14	.30	.19	.56	.18	.42	.31	.73
<i>p</i> = 3	supLM	.15	.48	.18	.52	.15	.48	.18	.52
	SupLM ^h	.15	.44	.18	.48	.15	.44	.18	.48
	AveLM	.08	.17	.19	.54	.08	.17	.19	.54
	AveLM ^h	.07	.16	.18	.47	.07	.16	.18	.47
	STAR	.10	.21	.15	.43	.10	.21	.15	.43

deterioration in power from the use of the heteroskedasticity-consistent test statistics, rather than the tests which assume homoskedasticity. The power loss is fairly mild, however. In sum, the simulation evidence strongly favors using our new tests.

6. APPLICATION TO U.S. GNP

We now apply this testing methodology to Potter's (1995) model of U.S. GNP. He used a SETAR in growth rates with a first, second, and fifth autoregressive lag. To select the threshold and delay parameters, he used informal graphical methods. While not an optimal estimation method, this is still conditional on the data, and hence invalidates the use of conventional test statistics and critical values. Our tests, on the other hand, allow a direct assessment of the statistical significance of his model.

The data is real GNP (seasonally adjusted) from Citibase for the period 1947-1990 and transformed into annualized quarterly growth rates. We estimated the model by least squares, allowing the threshold variable γ to vary from the 15th to the 85th percentile of the empirical distribution of x_t , and the delay parameter d over 1, 2, and 5. Our estimates are $\hat{d} = 2$, $\hat{\gamma} = 0.01$, and:

Regime 1— $x_{t-2} \leq 0.01$:

$$x_t = \underset{(1.12)}{-3.21} + \underset{(.25)}{.51} x_{t-1} - \underset{(.31)}{.93} x_{t-2} - \underset{(.25)}{.38} x_{t-5} + \hat{\epsilon}_t, \quad \hat{\sigma}_1^2 = 23.5;$$

TABLE IV
TESTS FOR THRESHOLD EFFECT IN U.S. GNP: 1947–1990

	SupLM	ExpLM	AveLM	SupLM ^h	ExpLM ^h	AveLM ^h
g_n	18.2	4.8	4.6	14.1	4.0	4.7
$\hat{\rho}_n^j$	0.04	0.09	0.29	0.17	0.17	0.27

Regime 2— $x_{t-2} > 0.01$:

$$x_t = 2.14 + .30 x_{t-1} + .18 x_{t-2} - .16 x_{t-5} + \hat{\varepsilon}_t, \quad \hat{\sigma}_2^2 = 12.1.$$

(.77) (.10) (.10) (.07)

Heteroskedastic-consistent standard errors are given in parenthesis.

The LM-based tests for the hypothesis of a single regime (no threshold effect) are reported in Table IV, with mixed results. If the homoskedasticity hypothesis is maintained, then supLM is marginally significant at the asymptotic 5% level, while aveLM is far from the rejection region. The point estimates for the error variance in the two regimes suggest that there may indeed be error heteroskedasticity. The test statistics which are robust to heteroskedasticity, however, are all far from standard rejection regions. Should we believe the “rejection” implied by supLM or the caution implied by the other statistics? The marginally significant supLM statistic could possibly be due to heteroskedasticity, yet the insignificant supLM^h statistic could possibly be due to its lower power. Our simulation evidence, however, suggested that the power loss is small. No definitive answer is possible, but it is fair to conclude that the apparent significance of the threshold model is consistent with sampling variation.

7. CONCLUSION

Many econometric models contain unidentified parameters under interesting parametric restrictions. Tests concerning these restrictions cannot use conventional statistical theory to assess significance. This paper extends the literature by developing a simple simulation technique which produces p -value statistics with an asymptotic uniform distribution under the null hypothesis.

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APPENDIX: MATHEMATICAL PROOFS

PROOF OF THEOREM 1: We start by showing that $S_n \Rightarrow S$. Note that for each $\gamma \in \Gamma$, $s_t(\gamma) = x_t(\gamma)\varepsilon_t$ is a square integrable stationary martingale difference, to which the pointwise central limit

theorem applies. Furthermore, note that the covariance kernel for $S_n(\gamma)$ is

$$E(S_n(\gamma_1)S_n(\gamma_2)') = \frac{1}{n} \sum_{t=1}^n E(s_t(\gamma_1)s_t(\gamma_2)') = K(\gamma_1, \gamma_2).$$

The multivariate central limit theorem establishes the finite dimensional distributional convergence.

To establish stochastic equicontinuity, we appeal to Theorem 1, Application 5, of Doukhan, Massart, and Rio (1994). Our data, and hence the summands $s_t(\gamma)$, satisfy the necessary absolute regularity mixing decay rate, and the envelope function is $\sup_\gamma |s_t(\gamma)| = \bar{x}_t |\varepsilon_t|$, which is L^{2v} bounded under Assumption 1. Finally, we need to show that the log of the L^{2v} bracketing numbers is integrable. Since $\Gamma \subset R^q$, one can always find a set Γ_N and constant $G < \infty$ so that for all γ there is some $\gamma_k \in \Gamma_N$ satisfying

$$(8) \quad |\gamma - \gamma_k| \leq GN^{-1/q}.$$

Set $N(\delta) = G^q B^q / \lambda^q \delta^{q/\lambda}$. Using Assumption 2 and (8), for all $\gamma \in \Gamma$,

$$(9) \quad \begin{aligned} \|s_t(\gamma) - s(\gamma_k)\|_{2v} &= \|(h_t(\gamma) - h_t(\gamma_k))\varepsilon_t\|_{2v} \\ &\leq B|\gamma - \gamma_k|^\lambda \leq BG^\lambda N^{-\lambda/q} = \delta. \end{aligned}$$

Thus $N(\delta)$ satisfy the definition of the L^{2v} bracketing numbers. Since

$$\int_0^1 \sqrt{\log(N(\delta))} d\delta \leq \sqrt{q \log(GB^{1/\lambda})} + \int_0^1 \sqrt{\log(\delta^{-q/\lambda})} d\delta < \infty$$

the conditions of Doukhan, Massart, and Rio are met, establishing that S_n is stochastically equicontinuous, and hence that $S_n \Rightarrow S$.

Standard algebra and Assumptions 1–3 show that

$$(10) \quad \begin{aligned} \sqrt{n} R' \hat{\beta}(\gamma) &= R' M_n(\gamma, \gamma)^{-1} S_n(\gamma) + R' M_n(\gamma, \gamma)^{-1} M_n(\gamma, \gamma_0) R c \\ &\Rightarrow R' M(\gamma, \gamma)^{-1} S(\gamma) + R' M(\gamma, \gamma)^{-1} M(\gamma, \gamma_0) R c = \bar{S}(\gamma) + \bar{Q}(\gamma) c. \end{aligned}$$

It is not hard to show that $\hat{V}_n(\gamma) \rightarrow K(\gamma, \gamma)$ a.s., uniformly over γ . Hence

$$(11) \quad R' \hat{V}_n^*(\gamma) R \rightarrow R' M(\gamma, \gamma)^{-1} K(\gamma, \gamma) M(\gamma, \gamma)^{-1} R = \bar{K}(\gamma, \gamma).$$

The continuous mapping theorem (CMT) applied to (10) and (11) show that $T_n \Rightarrow T^c$, and hence $g_n \Rightarrow g^c$ by the CMT. Q.E.D.

PROOF OF THEOREM 2: Let $s_t^*(\gamma) = s_t(\gamma)v_t$ and $S_n^*(\gamma) = (1/\sqrt{n})\sum_{t=1}^n s_t^*(\gamma)$. We first show that $S_n^* \Rightarrow_p S$. Let W denote the set of samples w for which

$$(12) \quad \limsup_{n \rightarrow \infty} \frac{1}{n} \sum_{t=1}^n \bar{x}_t^2 \varepsilon_t^2 < \infty,$$

$$K_n(\gamma_1, \gamma_2) \rightarrow K(\gamma_1, \gamma_2), \quad \text{uniformly over } \gamma_1 \text{ and } \gamma_2.$$

Under Assumptions 1–3, $P(W) = 1$. Take any $w \in W$. For the remainder of the proof, we will be operating conditionally on w , so all of the randomness appears in the iid $N(0, 1)$ variables v_t . Set $E_w X = E(X|w)$.

Note that $s_t^*(\gamma)$ are independent mean-zero normal random vectors, and S_n^* is a mean-zero Gaussian process with covariance function

$$E_w(S_n^*(\gamma_1)S_n^*(\gamma_2)') = \frac{1}{n} \sum_{t=1}^n E_w(s_t(\gamma_1)s_t(\gamma_2)'v_t^2)$$

$$= \frac{1}{n} \sum_{t=1}^n s_t(\gamma_1) s_t(\gamma_2)' = K_n(\gamma_1, \gamma_2).$$

Since $K_n(\gamma_1, \gamma_2) \rightarrow K(\gamma_1, \gamma_2)$ by (12), the finite dimensional distribution of S_n^* converge to those of S .

Next, the envelope of $s_t(\gamma)$ is L^2 -integrable:

$$\limsup_{n \rightarrow \infty} \frac{1}{n} \sum_{t=1}^n E_w \sup_{\gamma \in \Gamma} |s_t(\gamma)|^2 = \limsup_{n \rightarrow \infty} \frac{1}{n} \sum_{t=1}^n \bar{x}_t^2 \varepsilon_t^2 < \infty,$$

by (12). Define the norms

$$\begin{aligned} \rho_n(\gamma_1, \gamma_2) &= (E_w |S_n^*(\gamma_1) - S_n^*(\gamma_2)|^2)^{1/2} \\ &= \left(\frac{1}{n} \sum_{t=1}^n (s_t(\gamma_1) - s_t(\gamma_2))^2 \right)^{1/2} \end{aligned}$$

and $\rho(\gamma_1, \gamma_2) = \|s_t(\gamma_1) - s_t(\gamma_2)\|_2$. Observe that

$$\begin{aligned} \rho_n(\gamma_1, \gamma_2) &= \left[\text{tr} \left\{ \frac{1}{n} \sum_{t=1}^n (s_t(\gamma_1) - s_t(\gamma_2))(s_t(\gamma_1) - s_t(\gamma_2))' \right\} \right]^{1/2} \\ &= [\text{tr}\{K_n(\gamma_1, \gamma_1) + K_n(\gamma_2, \gamma_2) - K_n(\gamma_2, \gamma_1) - K_n(\gamma_1, \gamma_2)\}]^{1/2} \\ &\rightarrow [\text{tr}\{K(\gamma_1, \gamma_1) + K(\gamma_2, \gamma_2) - K(\gamma_2, \gamma_1) - K(\gamma_1, \gamma_2)\}]^{1/2} = \rho(\gamma_1, \gamma_2) \end{aligned}$$

uniformly over γ_1 and γ_2 under (12).

For any integer N , let Γ_N be the set satisfying (8), and for any $\delta > 0$ set $N(\delta) = G^q B^{q/\lambda} \delta^{q/\lambda}$. By the monotonicity of the L^p -norms and (9), we have

$$\rho(\gamma, \gamma_k) \leq \|s_t(\gamma) - s_t(\gamma_k)\|_{2\nu} \leq \delta,$$

establishing that $N(\delta)$ are the L^2 bracketing numbers, and have an exponential decay rate. Pollard (1990, Theorem 10.6) showed that S_n^* is stochastically equicontinuous under these conditions. Hence (for the particular w we selected), $S_n^* \Rightarrow S$. Since $P(W) = 1$, $S_n^* \Rightarrow S$, as desired.

Second, we wish to show that $\hat{S}_n \Rightarrow_p S$. Note that $\hat{S}_n(\gamma) = S_n^*(\gamma) + S_n^r(\gamma)$, where $S_n^r(\gamma) = (1/\sqrt{n}) \sum_{t=1}^n x_t(\gamma)(\hat{\varepsilon}_t(\gamma) - \varepsilon_t)v_t$. We have

$$(13) \quad \sup_{\gamma \in \Gamma} |S_n^r(\gamma)| \leq \sup_{\gamma \in \Gamma} \left| \frac{1}{n} \sum_{t=1}^n x_t(\gamma)x_t(\gamma)'v_t \right| \sup_{\gamma \in \Gamma} |\sqrt{n}(\hat{\beta}(\gamma) - \beta_0)|.$$

Using the same arguments as before, we can see that conditional on w , $(1/n) \sum_{t=1}^n x_t(\gamma)x_t(\gamma)'v_t \Rightarrow 0$ a.s. Thus (13) converges weakly in probability to zero, and $\hat{S}_n \Rightarrow_p S$.

This allows us to find under Assumptions 1-3 and the CMT that

$$\begin{aligned} \hat{T}_n(\gamma) &= \hat{S}_n(\gamma)' M_n(\gamma, \gamma)^{-1} R(R' \hat{V}_n^*(\gamma) R)^{-1} R' M_n(\gamma, \gamma)^{-1} \hat{S}_n(\gamma) \\ &\Rightarrow_p \bar{S}(\gamma)' \bar{K}(\gamma, \gamma)^{-1} \bar{S}(\gamma) = T^0(\gamma). \end{aligned}$$

Thus $\hat{g}_n = g(\hat{T}_n) \Rightarrow_p g(T^0) = g^0$, which implies that $\hat{F}_n(x) \rightarrow_p F^0(x)$, uniformly in x . We conclude that

$$\hat{p}_n = 1 - \hat{F}_n(g_n) = 1 - F^0(g_n) + o_p(1) = p_n + o_p(1),$$

as stated. The proof for \hat{p}_n is similar.

Q.E.D.

The proof of Theorem 3 relies on the following uniform strong law.

LEMMA 1: *If $\{w_t\}$ is strictly stationary and ergodic, $E|\phi(w_t)| < \infty$, and w_t has a continuous distribution, then*

$$(14) \quad \sup_{\gamma \in R} \left| \frac{1}{n} \sum_{t=1}^n \phi(w_t) \{w_t \leq \gamma\} - E(\phi(w_t) \{w_t \leq \gamma\}) \right| \rightarrow 0 \quad a.s.$$

PROOF OF LEMMA 1: Pollard (1984, Theorem II.2) established (14) when for each $\varepsilon > 0$ there exists a class of approximating functions $\{f_{\varepsilon,k}^u(w_t), f_{\varepsilon,k}^l(w_t) : k = 1, \dots, K_\varepsilon\}$, $K_\varepsilon < \infty$, which have the property that for each γ , there exists some k such that $f_{\varepsilon,k}^l(w_t) \leq \phi(w_t) \{w_t \leq \gamma\} \leq f_{\varepsilon,k}^u(w_t)$ and $E|f_{\varepsilon,k}^u(w_t) - f_{\varepsilon,k}^l(w_t)| < \varepsilon$. We can construct a qualifying set of approximating functions as follows. Set $K_\varepsilon = 2E|\phi(w_t)|/\varepsilon$ and select $\{\gamma_1, \dots, \gamma_{K_\varepsilon}\}$ so that for all k

$$(15) \quad E(|\phi(w_t)| \{\gamma_k < w_t \leq \gamma_{k+1}\}) \leq \varepsilon/2.$$

(This is possible since w_t has a continuous distribution.) Then set

$$f_{\varepsilon,k}^u(w_t) = \phi(w_t) \{w_t \leq \gamma_{k+1}\} \{ \phi(w_t) \geq 0 \} + \phi(w_t) \{w_t \leq \gamma_k\} \{ \phi(w_t) < 0 \}$$

and

$$f_{\varepsilon,k}^l(w_t) = \phi(w_t) \{w_t \leq \gamma_k\} \{ \phi(w_t) \geq 0 \} + \phi(w_t) \{w_t \leq \gamma_{k+1}\} \{ \phi(w_t) < 0 \}.$$

By construction, for all γ , there is some k such that $\gamma_k < \gamma \leq \gamma_{k+1}$ and $f_{\varepsilon,k}^l(w_t) \leq \phi(w_t) \{w_t \leq \gamma\} \leq f_{\varepsilon,k}^u(w_t)$. Hence,

$$\begin{aligned} E|f_{\varepsilon,k}^u(w_t) - f_{\varepsilon,k}^l(w_t)| &\leq 2E|\phi(w_t) (\{w_t \leq \gamma_{k+1}\} - \{w_t \leq \gamma_k\})| \\ &= 2E(|\phi(w_t)| \{\gamma_k < w_t \leq \gamma_{k+1}\}) \leq \varepsilon, \end{aligned}$$

where the final inequality is (15).

Q.E.D.

PROOF OF THEOREM 3: When $\gamma' < \gamma$, by Holder's inequality and the boundedness of f ,

$$\begin{aligned} \|(h_t(\gamma) - h_t(\gamma')) \varepsilon_t\|_{2\nu} &= \|x_t \varepsilon_t \{ \gamma' < z_t \leq \gamma \}\|_{2\nu} \\ &\leq \|x_t \varepsilon_t\|_{2r} \| \{ \gamma' < z_t \leq \gamma \} \|_{2r\nu/(r-\nu)} \\ &\leq \|x_t\|_{4r} \|\varepsilon_t\|_{4r} \left(\int_{\gamma'}^{\gamma} f(x) dx \right)^{(r-\nu)/2r\nu} \leq B|\gamma - \gamma'|^\lambda \end{aligned}$$

where $B = \|x_t\|_{4r} \|\varepsilon_t\|_{4r} \bar{f}^{(r-\nu)/2r\nu}$ and $\lambda = (r-\nu)/2r\nu$. The same inequality holds for $\gamma < \gamma'$, establishing Assumption 2.

Note that

$$M_n(\gamma_1, \gamma_2) = \left[\begin{array}{cc} \frac{1}{n} \sum_{t=1}^n x_t x_t' & \frac{1}{n} \sum_{t=1}^n x_t x_t' \{z_t \leq \gamma_2\} \\ \frac{1}{n} \sum_{t=1}^n x_t x_t' \{z_t \leq \gamma_1\} & \frac{1}{n} \sum_{t=1}^n x_t x_t' \{z_t \leq \min(\gamma_1, \gamma_2)\} \end{array} \right] \rightarrow M(\gamma_1, \gamma_2)$$

and

$$K_n(\gamma_1, \gamma_2) = \begin{bmatrix} \frac{1}{n} \sum_{t=1}^n x_t x_t' \varepsilon_t^2 & \frac{1}{n} \sum_{t=1}^n x_t x_t' \varepsilon_t^2 \{z_t \leq \gamma_2\} \\ \frac{1}{n} \sum_{t=1}^n x_t x_t' \varepsilon_t^2 \{z_t \leq \gamma_1\} & \frac{1}{n} \sum_{t=1}^n x_t x_t' \varepsilon_t^2 \{z_t \leq \min(\gamma_1, \gamma_2)\} \end{bmatrix} \rightarrow K(\gamma_1, \gamma_2)$$

a.s., uniformly in γ_1 and γ_2 by Lemma 1, and the assumption that z_t has a continuous distribution. This establishes Assumption 3. Q.E.D.

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