

The empirical process of autoregressive residuals

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Abstract: The asymptotic theory of the residual empirical process of autoregressions with an intercept is developed. In contrast to situations without intercept the asymptotic distribution does not depend on the location of the characteristic roots. This is important in applications, as the question of the distribution of the innovations then can be addressed without having to locate the characteristic roots. As a second contribution it is shown how the distribution of residual empirical moments are easily derived from the distribution of the residual empirical process.

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1 Introduction

The asymptotic theory of the empirical process of autoregressive residuals is revisited. The focus is on autoregressions including intercepts as in most applications. In contrast to situations without intercept it turns out that the asymptotic distribution does not depend on the location of the characteristic roots. This is important in applications, as the question of the distribution of the innovations then can be addressed without having to locate the characteristic roots. A second contribution is to exploit the simple, but apparently unnoticed, expression for moments as integrals with respect to the distribution function. This implies that the distribution of the empirical moments of the residuals can easily be derived from the distribution of the empirical process.

The simple first order autoregression without intercept has been studied extensively. Boldin (1981) and Koul and Leventhal (1989) studied the stationary and the explosive case respectively and found the same Gaussian limiting distribution as in regressions with fixed regressors but no intercept. Ling (1998) and Lei and Wei (1999) studied the case with a unit root and a known scale and found the limiting distribution to be non-Gaussian in that case. Thus, for the simple first order autoregression it is not possible to separate the questions of the distribution of the characteristic roots and of the location of the characteristic roots.

The autoregression with intercept has received less attention. Pierce (1985) and Koul and Leventhal (1989) studied the stationary and the explosive case respectively and found the same Gaussian limiting distribution as in regressions with fixed regressors including an intercept. The unit root case has not been studied. In this paper it is shown that the limiting distribution is the same,

so that the limiting distribution is invariant to the location of the characteristic roots. The key to the result is that the non-Gaussian component in the case of no intercept stems from the sum of the residuals. When an intercept is included in the model, as in most applications, this sum is zero and the non-Gaussian component does not arise. A similar parameter invariance apply for the likelihood-based tests for unit roots and for the order of autoregressions, but not for the usual correlograms, see Nielsen (2001, 2006a,b).

The proofs presented here draw on the results of Lee and Wei (1999) and Koul (2002) for residual empirical processes for stochastic regressions. Some work is needed since neither include an intercept and the former has an known scale. By combining this work with the asymptotic theory of autoregressive estimators developed in Nielsen (2005) it is possible to consider autoregressive distributed lag models with general deterministic terms, which are frequently used in practice.

The simple autoregression has received some further attention in the stationary case with no intercept and known scale. In that situation the limiting process is a Brownian bridge. The partial empirical process has been proved to converge to a Kiefer process by Bai (1994) and recently by Na, Lee and Park (2005) for the case with measurement errors. Since the empirical process converges to a Brownian bridge the Kolmogorov-Smirnov test is distribution free. Therefore, the empirical process can be used not only for testing for a given reference distribution of the innovations but also to estimate the innovation distribution. When the scale is unknown the Kolmogorov-Smirnov statistic is not distribution free, but a distribution free statistic can be found using the Khmaladze transformation involving the innovation density. If the focus is to estimate the innovation distribution this density will also have to be estimated, which is challenging unless the sample size is large. Bai (2003) points out that the Khmaladze transformation apply for autoregressions, but focuses on the testing issue for non-linear time series where the transformation can simplify complicated limiting distributions.

The considered class of autoregressions is quite general. Autoregressive distributed lag models are allowed where the joint distribution of the included variables is vector autoregressive without restrictions on the characteristic roots. The deterministic components can be polynomial and seasonal. The main result states the asymptotic properties of the residual empirical process. As a corollary the empirical moments are analysed. Focusing on the Gaussian case this is used to show how the innovation distribution can be tested using a Kolmogorov-Smirnov test, an Anderson-Darling-type test, a Probability-Probability plot or a Quantile-Quantile plot with test bands, or the standard cumulant based tests for normality, known as the Jarque-Bera tests. Finally, an empirical illustration follows.

2 The empirical process

A general autoregressive model is set up. The asymptotic theory of the empirical distribution of the residuals then follows.

2.1 The general autoregressive model

Let X_t be a p -dimensional time series partitioned in terms of a univariate time series Y_t and $(p - 1)$ -dimensional time series Z_t . The general univariate model is given by

$$Y_t = \rho Z_t + \sum_{j=1}^k \alpha_j Y_{t-j} + \sum_{j=1}^k \beta_j' Z_{t-j} + \nu D_{t-1} + \sigma \varepsilon_t, \quad (t = 1, \dots, T) \quad (2.1)$$

conditional on X_0, \dots, X_{1-k} , with independent innovations ε_t with distribution function F . The term D_{t-1} is a deterministic term, which will be discussed in further detail below. When ρ is restricted to zero, so Z_t is absent on the right hand side, this is the marginal equation of a vector autoregression, and when ρ is unrestricted this is an autoregressive distributed lags model. If the Z process is omitted this reduces to an autoregression. Note, that when $k = 0$ and $D_{t-1} = 1$ the model reduces to a classical regression model with an intercept. In addition, when ρ is restricted to zero, this becomes a location-scale model for Y_t .

Least squares estimation of the equation (2.1) gives the scaled residuals

$$\hat{\varepsilon}_t = \frac{Y_t - \hat{\rho} Z_t - \sum_{j=1}^k \hat{\alpha}_j Y_{t-j} - \sum_{j=1}^k \hat{\beta}_j' Z_{t-j} - \hat{\nu} D_{t-1}}{\hat{\sigma}},$$

if, for instance, ρ is unrestricted. The empirical distribution function of the residuals is defined as

$$\hat{F}(x) = \frac{1}{T} \sum_{t=1}^T 1_{(\hat{\varepsilon}_t \leq x)},$$

with the associated empirical process

$$\hat{\mathbb{F}}\{F(x)\} = \sqrt{T} \left\{ \hat{F}(x) - F(x) \right\}, \quad (2.2)$$

which has argument $u = F(x)$ on the unit interval.

The inclusion of the intercept plays a crucial role in the analysis of the empirical process. The issue is that the sub-sequent asymptotic expansion of $\hat{\mathbb{F}}$ involves the term $T^{-1/2} \sum_{t=1}^T \hat{\varepsilon}_t$. When no intercept is included this can be shown to vanish as long as the characteristic roots are different from one, whereas manipulations as in Chan and Wei (1988) shows that it has a Dickey-Fuller-type distribution in the presence of unit roots. When an intercept is included the likelihood equation for the intercept implies that $\sum_{t=1}^T \hat{\varepsilon}_t = 0$ regardless of the values of the autoregressive parameter.

In order to discuss the distribution of the empirical process the joint distribution of the time series $X_t = (Y_t, Z_t)'$ has to be specified. If this is assumed to satisfy a vector autoregression the results for general vector autoregressions given by Nielsen (2005) can be used. That paper is a generalisation of the work by Lai and Wei (1985), who did not consider deterministic terms. Thus, suppose the time series X_t and the deterministic series D_t satisfy the autoregressive equations

$$X_t = \sum_{j=1}^k A_j X_{t-j} + \mu D_{t-1} + \xi_t, \quad (t = 1, \dots, T), \quad (2.3)$$

$$D_t = \mathbf{D} D_{t-1}, \quad (2.4)$$

conditional on X_0, \dots, X_{1-k} , where the vector innovations, ξ_t , are partitioned as $\xi_t = (\xi_{y,t}, \xi'_{z,t})'$. The vector innovations are assumed to have mean zero and a positive definite variance matrix partitioned as

$$\text{Var}(\xi_t) = \Omega = \begin{pmatrix} \Omega_{yy} & \Omega_{yz} \\ \Omega_{zy} & \Omega_{zz} \end{pmatrix}. \quad (2.5)$$

The univariate model (2.1) can then arise in two ways:

1. When ρ is restricted to zero then (2.1) is simply the first equation in (2.3) so $\sigma\varepsilon_t = \xi_{y,t}$ with mean zero and variance $\sigma^2 = \Omega_{yy}$.
2. When $\rho = \Omega_{yz}\Omega_{zz}^{-1}$ and ξ is normally distributed, then (2.1) states the conditional model for Y_t given Z_t and the past, with $\sigma\varepsilon_t = \xi_{y,t} - \rho\xi_{z,t}$ with mean zero and variance $\sigma^2 = \Omega_{yy} - \Omega_{yz}\Omega_{zz}^{-1}\Omega_{zy}$.

The formulation for the deterministic term D_t allows a joint autoregressive companion representation of X_t, D_t , and is inspired by Johansen (2000). The matrix \mathbf{D} has characteristic roots on the complex unit circle, so D_t is a vector of terms such as a constant, a linear trend, or periodic functions like seasonal dummies. For example,

$$\mathbf{D} = \begin{pmatrix} 1 & 0 \\ 1 & -1 \end{pmatrix} \quad \text{with} \quad D_0 = \begin{pmatrix} 1 \\ 1 \end{pmatrix}$$

will generate a constant and a dummy for a bi-annual frequency. The deterministic term D_t is assumed to have linearly independent coordinates, which is formalised as follows.

Assumption 2.1 $|\text{eigen}(\mathbf{D})| = 1$ and $\text{rank}(D_1, \dots, D_{\dim \mathbf{D}}) = \dim \mathbf{D}$.

To ensure that an intercept is included in the model an additional assumption to \mathbf{D} is needed.

Assumption 2.2 Assume \mathbf{D} has at least one eigenvalue of unity.

Note that indicator dummies for single observations of the form $D_t = 1_{(t=j)}$ can also be included in the analysis. These will result in a zero residual at observations j , that is $\widehat{\varepsilon}_j = 0$. Such residuals are simply omitted when forming the empirical distribution function.

2.2 Asymptotic theory for the empirical process

To formulate the main result some assumptions to innovations ξ_t and ε_t are needed. The innovations, ξ_t , of the vector autoregression (2.3) are assumed to satisfy a martingale difference sequence assumption to exploit the consistency results of Nielsen (2005).

Assumption 2.3 Suppose the sequence of innovations ξ_t of the vector autoregression (2.3) is a martingale difference sequence with respect to an increasing filtration \mathcal{F}_t so $\mathbf{E}(\xi_t|\mathcal{F}_{t-1}) = 0$ and where the initial values X_0, \dots, X_{1-k} are \mathcal{F}_0 -measurable and

$$\sup_t \mathbf{E}\{(\xi_t'\xi_t)^{\lambda/2}|\mathcal{F}_{t-1}\} \stackrel{a.s.}{<} \infty \quad \text{for some } \lambda > 4, \quad (2.6)$$

$$\mathbf{E}(\xi_t\xi_t'|\mathcal{F}_{t-1}) \stackrel{a.s.}{=} \Omega \quad \text{where } \Omega \text{ is positive definite.} \quad (2.7)$$

Assumption 2.3 exclude the possibility that the innovations could be autoregressive conditional heteroscedastic (ARCH). This assumption is primarily needed to cover the explosive case. It is therefore necessary for the coveted parameter invariance in the main result.

A further set of three assumptions are needed for the innovations ε_t of the regression equation (2.1). It is not sufficient that these innovations are martingale difference sequences, essentially because the empirical process describes the entire distribution, and thus all moments of ε_t . The first assumption ensures the convergence of the uniform empirical process.

Assumption 2.4 Suppose the innovations ε_t of the regression equation (2.1) are independent and identically distributed so $\mathbf{E}\varepsilon_t = 0$, $\mathbf{Var}\varepsilon_t = 1$, and with marginal distribution function F .

A second assumption is a martingale assumption, which is needed for the case where ρ is unrestricted so the regression equation (2.1) becomes an autoregressive distributed lags equation.

Assumption 2.5 If ρ is unrestricted define the filtration \mathcal{G}_{t-1} as the sigma field over \mathcal{F}_{t-1} and Z_t , otherwise, if ρ is restricted to zero let $\mathcal{G}_{t-1} = \mathcal{F}_{t-1}$. Suppose the innovations ε_t are independent of \mathcal{G}_{t-1} .

The third assumption concerns the distribution function F of the innovations, ε_t .

Assumption 2.6 Suppose the distribution function F has density f which is positive and differentiable everywhere, and satisfies

$$\sup_{x \in \mathbf{R}} |xf(x)| < \infty, \quad \sup_{x \in \mathbf{R}} |f'(x)| < \infty, \quad \sup_{x \in \mathbf{R}} |x^2f'(x)| < \infty.$$

The Assumption 2.6 is satisfied by many distributions, notably the standard normal distributions. Denoting the standard normal density by φ it holds that $\varphi'(x) = -x\varphi(x)$. Since φ has exponentially declining tails the boundedness follows.

The Brownian bridge \mathbb{U} can be written in terms of a standard Brownian motion \mathbb{B} as $\mathbb{U}(u) = \mathbb{B}(u) - u\mathbb{B}(1)$. The asymptotic result for the empirical process $\widehat{\mathbb{F}}(u)$ is a generalisation of this result, where the limiting distribution is expressed in terms of stochastic integrals with respect to the Brownian bridge \mathbb{U} .

The limiting distribution of the empirical process will be expressed in terms of a stochastic integrals with respect to a Brownian bridges as discussed by Shorack and Wellner (1986, p. 91-95). It is worth noting that stochastic integrals $\int_0^1 h_j(u) d\mathbb{U}(u)$, are well-defined for square integrable functions h_1, h_2 . They are normally distributed with expectation zero, and variance given by

$$\begin{aligned} & \text{Cov} \left\{ \int_0^1 h_1(u) d\mathbb{U}(u), \int_0^1 h_2(u) d\mathbb{U}(u) \right\} \\ &= \int_0^1 h_1(u) h_2(u) du - \left\{ \int_0^1 h_1(u) du \right\} \left\{ \int_0^1 h_2(u) du \right\}. \end{aligned} \quad (2.8)$$

The Brownian bridge itself can be written as $\mathbb{U}(v) = \int_0^1 1_{(u \leq v)} d\mathbb{U}(u)$.

The main result can now be formulated. This shows that as long as an intercept is included in the autoregression the same Gaussian limit distribution applies for all values of the characteristic roots. Thus, it is possible to make inference about the distribution of the autoregressive innovations without knowing the location of the characteristic roots.

Theorem 2.7 *Suppose the model (2.3) with no restrictions to the characteristic roots and the Assumptions 2.1, 2.2, 2.3, 2.4, 2.5, 2.6 are satisfied. Then*

$$\widehat{\mathbb{F}}(\cdot) \xrightarrow{D} \mathbb{X}_{\mathbb{F}}(\cdot),$$

on $D[0, 1]$, where $\mathbb{X}_{\mathbb{F}}$ is the Gaussian process defined by

$$\mathbb{X}_{\mathbb{F}}(u) = \left[\begin{array}{c} 1 \\ \mathbf{f} \{ \mathbb{F}^{-1}(u) \} \\ \frac{1}{2} \mathbb{F}^{-1}(u) \mathbf{f} \{ \mathbb{F}^{-1}(u) \} \end{array} \right]' \int_0^1 \left[\begin{array}{c} 1_{[s \leq u]} \\ \mathbb{F}^{-1}(s) \\ \{ \mathbb{F}^{-1}(s) \}^2 \end{array} \right] d\mathbb{U}(s). \quad (2.9)$$

The process $\mathbb{X}_{\mathbb{F}}$ defined in (2.9) is the sum of three components. The first component is simply the Brownian Bridge $\mathbb{U}(u)$ arising from the uniform empirical process $\widehat{\mathbb{U}}$. The second and the third components relate to $\widehat{\mathbb{F}} - \widehat{\mathbb{U}}$ and arise due to the estimation of the variance and the intercept, respectively. The process $\mathbb{X}_{\mathbb{F}}$ is standard, in the sense of applying to cross-sectional regression problems including a constant, see Shorack & Wellner (1986, p.197f).

For a known reference distribution the process $\mathbb{X}_{\mathbb{F}}$ can be simulated in various ways. The first approach is to use the covariance structure of $\mathbb{X}_{\mathbb{F}}$. For any choice of reference distribution \mathbb{F} this covariance structure can be found using the formula (2.8). For the Gaussian case it is found in Theorem 2.9 below. For a particular grid over the interval $[0, 1]$, the covariance matrix is computed and its square root multiplied onto a vector of generated independent standard normal variables. The second approach is to compute the stochastic integrals in (2.9) directly. First, a Brownian motion \mathbb{B} is computed for a grid over $[0, 1]$ by taking partial sums of generated independent standard normal variables and is transformed into a Brownian bridge using the formula $\mathbb{U}(u) = \mathbb{B}(u) - u\mathbb{B}(1)$. Next, the integral (2.9) is formed. This approach is a little more convoluted, but numerically faster when dealing with a fine grid over $[0, 1]$.

2.3 Empirical moments

The asymptotic properties of the moments of the residuals can be found rather easily from the empirical process. This point seems to have escaped the literature. It will later, in §3.4, be used to establish the asymptotic behaviour of the usual cumulant-based tests for normality. Some notation is needed. Define the empirical moments of the standardised residuals,

$$\hat{\mu}_m = \frac{1}{T} \sum_{t=1}^T \hat{\varepsilon}_t^m = \int_{\mathbf{R}} x^m d\hat{\mathbf{F}}(x),$$

and the least squares estimator for the variance parameter, normalised by T ,

$$\hat{\sigma}^2 = \frac{1}{T} \sum_{t=1}^T \left(Y_t - \hat{\rho}Z_t - \sum_{j=1}^k \hat{\alpha}_j Y_{t-j} - \sum_{j=1}^k \hat{\beta}'_j Z_{t-j} - \hat{\nu}D_{t-1} \right)^2.$$

It then holds that $\hat{\mu}_2 = 1$, and, when including an intercept in the model, also $\hat{\mu}_1 = 0$. Defining the population moments

$$\mu_m = \int_{\mathbf{R}} x^m d\mathbf{F}(x) = \int_0^1 \{\mathbf{F}^{-1}(u)\}^m du, \quad (2.10)$$

the sample and population moments can be brought together as

$$\begin{aligned} \sqrt{T}(\hat{\mu}_m - \mu_m) &= \sqrt{T} \left\{ \int_{\mathbf{R}} x^m d\hat{\mathbf{F}}(x) - \int_{\mathbf{R}} x^m d\mathbf{F}(x) \right\} \\ &= \int_{\mathbf{R}} x^m d\hat{\mathbb{F}}_T(x) = \int_0^1 \{\mathbf{F}^{-1}(u)\}^m d\hat{\mathbb{F}}_T\{\mathbf{F}^{-1}(u)\}. \end{aligned} \quad (2.11)$$

The asymptotic theory then follows by replacing the empirical process $\hat{\mathbb{F}}_T$ with the limiting Gaussian process $\mathbb{X}_{\mathbf{F}}$ established in Theorem 2.7. To describe the resulting limiting normal distribution let

$$\begin{aligned} h_2(u) &= \{\mathbf{F}^{-1}(u)\}^2, \\ h_m(u) &= \{\mathbf{F}^{-1}(u)\}^m - m\mu_{m-1}\{\mathbf{F}^{-1}(u)\} - \frac{m}{2}\mu_m\{\mathbf{F}^{-1}(u)\}^2. \end{aligned}$$

Theorem 2.8 *Suppose the model (2.3) with no restrictions to the characteristic roots and the Assumptions 2.1, 2.2, 2.3, 2.4, 2.5, 2.6 are satisfied, and that \mathbf{F} has moments of order at least $2M$. Then*

$$\sqrt{T}(\hat{\sigma}^2 - \sigma^2, \hat{\mu}_3 - \mu_3, \dots, \hat{\mu}_M - \mu_M)$$

converge jointly in distribution to a normal distribution with mean zero. The limiting distribution can be represented as

$$\sqrt{T} \left(\frac{\hat{\sigma}^2}{\sigma^2} - 1 \right) \xrightarrow{D} \int_0^1 h_2(u) d\mathbb{U}(u) = \int_0^1 \{\mathbf{F}^{-1}(u)\}^2 d\mathbb{U}(u), \quad (2.12)$$

$$\sqrt{T}(\hat{\mu}_m - \mu_m) \xrightarrow{D} \int_0^1 h_m(u) d\mathbb{U}(u) = \int_0^1 \{\mathbf{F}^{-1}(u)\}^m d\mathbb{X}_{\mathbf{F}}(u), \quad (2.13)$$

and the covariance matrix has entries

$$\omega_{mn} = \text{Cov} \left\{ \int_0^1 h_m(u) d\mathbb{U}(u), \int_0^1 h_n(u) d\mathbb{U}(u) \right\}$$

which satisfy, for $m, n \geq 3$,

$$\begin{aligned} \omega_{22} &= \mu_4 - 1, & \omega_{2m} &= \mu_{m+2} - m\mu_3\mu_{m-1} - \mu_m \left\{ 1 + \frac{m}{2}(\mu_4 - 1) \right\}, \\ \omega_{mn} &= \mu_{m+n} - \mu_{m+1}\mu_{n+1} - \left(\frac{mn}{2} + 1 \right) \mu_m\mu_n + \frac{mn}{4}(\mu_4 - 3)\mu_m\mu_n \\ &\quad + (\mu_{m+1} - m\mu_{m-1})(\mu_{n+1} - n\mu_{n-1}) + \frac{mn}{2}\mu_3(\mu_{m-1}\mu_n + \mu_m\mu_{n-1}) \\ &\quad - \frac{n}{2}\mu_n \{ \mu_{m+2} - (m+1)\mu_m \} - \frac{m}{2}\mu_m \{ \mu_{n+2} - (n+1)\mu_n \}. \end{aligned}$$

The result can form the basis for a cumulant-based test of the reference distribution F in a location-scale model. For instance, if ε_t follows a standard Gaussian distribution or a t -distribution with known degrees of freedom, then $Y_t = \mu + \sigma\varepsilon_t$ is a location-scale model. The Gaussian case is studied in further detail in §2.4.

2.4 The Gaussian case

The case where the reference distribution F is standard Gaussian, denoted Φ is of special interest. At first, the covariance structure of the limiting Gaussian process \mathbb{X}_F is found. Subsequently, the moments of the empirical distribution are described.

Theorem 2.9 *Suppose $F(u) = \Phi(u)$ is the standard normal distribution function with density φ . Then $\mathbb{X}_F = \mathbb{X}_\Phi$ has covariance given by*

$$\text{Cov} \{ \mathbb{X}_\Phi(u), \mathbb{X}_\Phi(v) \} = u(1-v) - \varphi \{ \Phi^{-1}(u) \} \varphi \{ \Phi^{-1}(v) \} \left\{ \frac{\Phi^{-1}(u)\Phi^{-1}(v)}{2} + 1 \right\},$$

for $0 \leq u \leq v \leq 1$, or equivalently, for $x, y \in \mathbf{R}$ so $x \leq y$,

$$\text{Cov} [\mathbb{X}_\Phi \{ \Phi(x) \}, \mathbb{X}_\Phi \{ \Phi(y) \}] = \Phi(x) \{ 1 - \Phi(y) \} - \varphi(x) \varphi(y) \left(\frac{xy}{2} + 1 \right).$$

Theorem 2.8 presented the asymptotic distribution of the sample moments of the standardised residuals. This result simplifies considerably in the Gaussian case. In that case the population moments are

$$\mu_m = \int_{\mathbf{R}} x^m d\Phi(x) = \int_0^1 \{ \Phi^{-1}(u) \}^m du = \begin{cases} 0 & \text{for } m \text{ odd,} \\ (m-1)!! & \text{for } m \text{ even,} \end{cases} \quad (2.14)$$

where $(m-1)!! = (m-1)(m-3)\cdots 3 \cdot 1$ is the odd factorial. In particular $\mu_3 = 0$, $\mu_4 = 3$, and $\mu_{m+1} = m\mu_{m-1}$. This immediately gives the following result, noting that Assumptions 2.4 and 2.6 are satisfied in the normal case.

Corollary 2.10 *Suppose $F(u) = \Phi(u)$ is the standard normal distribution function. Then the asymptotic covariance matrix in Theorem 2.8 has entries, for $m, n \geq 3$,*

$$\omega_{22} = 2, \quad \omega_{2m} = 0, \quad \omega_{mn} = \mu_{m+n} - \mu_{m+1}\mu_{n+1} - \left(\frac{mn}{2} + 1\right) \mu_m \mu_n.$$

The property that $\omega_{2m} = 0$ shows that $\hat{\sigma}^2$ is asymptotically independent of all higher moments of the standardised residuals, which in turn implies that it is independent of the higher cumulants. This relates to the exact independence found by Fisher (1930) for samples of independent normal variates. Note also that $\omega_{mn} = 0$ when m is even and n is odd, implying that for instance the third and fourth cumulants of the residuals are asymptotically independent.

3 Applications

Theorem 2.7 has a wide range of applications that can be helpful in autoregressive analysis. In the following, the empirical process is at first transformed into a standardised empirical process and a quantile process. Next, Kolmogorov-Smirnov and Anderson-Darling tests are discussed. Then test bands are presented for Probability-Probability and Quantile-Quantile plots. These are frequently used in econometrics, but without test bands. Finally, the frequently used cumulant based tests for normality are discussed.

3.1 The standardised empirical process and the quantile process

In Theorem 2.7, a Gaussian approximation was found for the distribution of the empirical process. Using this result, a standardised version of the empirical process can be defined

$$\hat{Z}_F(u) = \frac{\hat{F}(u)}{\widehat{\text{std}}\{\mathbb{X}_F(u)\}} \quad \text{for } u \in (0, 1).$$

This process is not defined at the end points $u = 0$ and $u = 1$ since both the empirical process and its variance are then zero. For each u in the open interval $(0, 1)$ Theorem 2.7 implies the pointwise convergence result:

$$\hat{Z}_F(u) \xrightarrow{D} Z_F(u) = \frac{\mathbb{X}_F(u)}{\text{std}\{\mathbb{X}_F(u)\}}. \quad (3.1)$$

The sequence of standardised empirical processes, \hat{Z}_F , is, however, not tight. The issue is the variation near 0 and near 1. In this way, Čibisov (1966) showed that

$$\sup_{0 < u < 1} \left| \frac{\hat{U}(u)}{\text{std}\{\mathbb{U}(u)\}} \right| \xrightarrow{P} \infty.$$

The rate of divergence has been studied intensively, with an overview given in Shorack and Wellner (1986). The process \hat{Z}_F will, however, converge on $D[b, c]$ where $[b, c]$ is an arbitrary closed subinterval of $(0, 1)$.

Rather than looking at the distribution function for the residuals, it is often of interest to look at their quantiles, defined in terms of the inverse distribution function, \widehat{F}^{-1} . The standardised quantile process is defined as

$$\widehat{Q}_F(x) = f(x) \sqrt{T} \left[\widehat{F}^{-1} \{F(x)\} - x \right]. \quad (3.2)$$

Using the functional δ -method, in an argument made more formally by Shorack and Wellner (1986, Chapter 18), it holds that

$$\widehat{Q}_F(x) \xrightarrow{D} \mathbb{X}_F(x), \quad (3.3)$$

on $D[b, c]$ where $[b, c] \subset \mathbf{R}$. The quantile process itself then satisfies

$$\widehat{R}_F(x) = \sqrt{T} \left[\widehat{F}^{-1} \{F(x)\} - x \right] \xrightarrow{D} \frac{\mathbb{X}_F(x)}{f(x)}. \quad (3.4)$$

on $D[b, c]$. In particular it holds that the standardised quantile process and the standardised empirical process have the same limiting distribution.

3.2 Test statistics based on the empirical process

In many applications it is convenient to summarize the empirical process in a single statistic. There are two reasons for this. *First*, for the uniform empirical process, distribution free tests can be constructed based on, for instance, the Kolmogorov-Smirnov statistic. This does, however, not seem possible when the location and scale are unknown as here. *Secondly*, tests can be constructed by evaluating the empirical process at a single point. Looking at two different points, x_1 and x_2 , the test statistics will be dependent. To achieve a correct size this has to be taken into account. This is done by looking at summary statistics such as the Kolmogorov-Smirnov statistic. Different summary statistics are found depending on the choice of weight function that is applied to the empirical process over the points $x \in \mathbf{R}$. There is no optimal way of choosing a weight function. Likewise, such statistics are bound to be dependent, so when using several statistics it is not clear how to achieve a correct size.

Kolmogorov-Smirnov-type statistics are formed from the empirical process as

$$\widehat{D}_F = \sup_{x \in \mathbf{R}} \left| \widehat{F} \{F(x)\} \right| = \sup_{0 \leq u \leq 1} \left| \widehat{F}(u) \right| \xrightarrow{D} \sup_{0 \leq u \leq 1} |\mathbb{X}_F(u)| = D_F, \quad (3.5)$$

$$\widehat{D}_F^+ = \sup_{x \in \mathbf{R}} \widehat{F} \{F(x)\} = \sup_{0 \leq u \leq 1} \widehat{F}(u) \xrightarrow{D} \sup_{0 \leq u \leq 1} \mathbb{X}_F(u) = D_F^+, \quad (3.6)$$

where the limiting distributions arise by applying the Continuous Mapping Theorem to the asymptotic distribution for the empirical process reported in Theorem 2.7. These statistics put most weight on values of u closest to 0.5. The limiting distributions of D_F , D_F^+ depend on the distribution of F . This is in contrast to the situation where the empirical distribution of independent, identically variables is compared to a known distribution. In that situation only the uniform term \widehat{U} is of relevance and a time transformation argument can be employed.

	E	Var	E log	Var log	50%	80%	90%	95%	97.5%	99%
D_{Φ}	0.631	0.0219	-0.487	0.0526	0.612	0.748	0.830	0.903	0.971	1.053
D_{Φ}^+	0.547	0.0231	-0.640	0.0731	0.524	0.666	0.753	0.831	0.903	0.990

Table 1: Simulated distribution of \widehat{D}_{Φ} and \widehat{D}_{Φ}^+

When testing for normality the distributions D_{Φ} and D_{Φ}^+ are of relevance. These were previously tabulated by Stephens (1974, Table 1A, case 3). Stephens' experiment was repeated using a fine grid with 10^4 points and 10^6 repetitions giving the numbers reported in Table 1. A convenient approximation to the p-values can be found using a Gamma distribution with the reported mean and variance as done for instance for a Dickey-Fuller F-type distribution in Nielsen (1997). Due to the extreme value nature of the Kolmogorov-Smirnov statistic, the Gamma approximation will under-fit the extreme upper tail slightly. For the reported quantiles the relative error of the Gamma quantiles compared to the simulated quantiles was at most 3.0%, and at most 1.8% when excluding the two most extreme quantiles. It was also attempted to fit a Weibull distribution, as the asymptotic distribution of the D_{Φ}^+ statistic based on the uniform empirical process is Weibull, see Billingsley (1968, p.85). The Weibull distribution can be fitted using mean and variance of the log-transformed statistic, see Johnson, Kotz, and Balakrishnan (1994, §21.4). For the empirical process of residuals the fit is, however, much worse than the Gamma fit, with relative errors of up to 17%.

While the Kolmogorov-Smirnov statistic puts most weight on deviations in the middle of the distribution, Anderson and Darling (1952) considered the possibility of constructing Kolmogorov-Smirnov-type statistics for the standardised empirical process, $\widehat{Z}_{\mathbb{F}}$, see (3.1). Anticipating the result of Čibisov (1966), they suggested taking supremum over a closed interval in the interior of the unit interval. Considering a symmetric interval for simplicity, this gives

$$\widehat{K}_{\mathbb{F},a} = \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} \left| \widehat{Z}_{\mathbb{F}}(u) \right| \xrightarrow{\text{D}} \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} |Z_{\mathbb{F}}(u)| = K_{\mathbb{F},a}.$$

A Kolmogorov-Smirnov-type statistics can be constructed in the same way for the empirical quantile process

$$\widehat{R}_{\mathbb{F},a} = \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} \left| \widehat{\mathbb{R}}_{\mathbb{F}}(u) \right| \xrightarrow{\text{D}} \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} |\mathbb{R}_{\mathbb{F}}(u)| = R_{\mathbb{F},a},$$

whereas the Anderson-Darling-type statistic for the empirical quantile process satisfies

$$\widehat{Q}_{\mathbb{F},a} = \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} \left| \frac{\widehat{\mathbb{R}}_{\mathbb{F}}(u)}{\widehat{\text{std}} \left\{ \widehat{\mathbb{R}}_{\mathbb{F}}(u) \right\}} \right| \xrightarrow{\text{D}} \sup_{\frac{1}{2}-a < u < \frac{1}{2}+a} |Z_{\mathbb{F}}(u)| = K_{\mathbb{F},a},$$

and has the same limiting distribution as the Kolmogorov-Smirnov statistic.

When testing for normality the distributions $K_{\Phi,a}$ and $R_{\Phi,a}$ are of relevance. These were simulated for a range of a values, a fine grid over u with 10^4 points

		1	a	a^2	a^{-1}	a^{-2}	$(\frac{1}{2} - a)^{-1}$	$(\frac{1}{2} - a)^{-2}$
$K_{\Phi,a}$	E	1.969	1.405		-0.04406	0.0008920	0.00028	
	Var	0.365	-0.291	0.141			-0.00001	
$R_{\Phi,a}$	E	1.672	-1.583	4.904	-0.02773		0.03722	-0.0000677
	Var	0.135	0.053	-1.293			0.02743	-0.0000151

Table 2: Response surface in a for expectation and variance of $\widehat{K}_{\Phi,a}$ and $\widehat{R}_{\Phi,a}$.

and 10^6 repetitions. Response surfaces in a for the expectation and variance of $K_{\Phi,a}$ and $R_{\Phi,a}$ are reported in Table 2 using 22 values of a chosen as $(0.05:0.40, 0.05)$, $(0.41:0.49, 0.01)$, $(0.491:0.495, 0.001)$. In all cases, the R^2 of the fits exceed 0.9995. It is not advisable to extrapolate the response surface for values of a outside $0.01 < a < 0.495$. For a given value of a the expectation and variance are computed and the distribution approximated using a Gamma distribution as above. It is evident that the response surfaces diverge for $a \rightarrow 0.5$. Different choices of a will emphasize different departures from normality. Due to the Čibisov result the literature does not give any guidance towards the choice of a as a function of the sample size.

3.3 P-P and Q-Q plots

At present, probability-probability and quantile-quantile plots, also called P-P and Q-Q plots, are used without an indication of the uncertainty. Based on the work present here, two types of test bands can be derived: pointwise bands and simultaneous bands based on the statistics presented in §3.2. Pointwise bands would be used in situations where the nature of the departure from the reference distribution is unknown, whereas the simultaneous bands are used to detect more specific types of departures.

Probability-probability plots are plots of $\widehat{u} = \widehat{F}(x)$ against $u = F(x)$ on a $[0, 1] \times [0, 1]$ -square. Test bands of the type $u \pm n^{-1/2}c_\alpha\sigma_u$ can be constructed using the results from above. Three different bands are worth noting.

Pointwise test bands can be established from Theorem 2.7, where c_α is the $1 - \alpha/2$ quantile of the standard normal distribution and σ_u^2 is the variance $\text{Var}\{\mathbb{X}_F(u)\}$ at the point u . For the standard normal case, $F = \Phi$, the variance is reported in Theorem 2.9. For other reference distributions it can be computed using (2.8).

Kolmogorov-Smirnov bands can be constructed by $\sigma_u = 1$ and choosing c_α from limiting distributions of the Kolmogorov-Smirnov statistics. This gives bands that are straight lines parallel to the 45°-line. This emphasises departures in the middle of the distribution. Bands of “two-sided” nature are found by choosing c_α as the $1 - \alpha$ quantile of D_F , whereas a “one-sided” are found from the $1 - \alpha/2$ quantile of D_F^+ .

Anderson-Darling bands are constructed by first choosing a value of a (perhaps 0.3, 0.4, or 0.45). Then σ_u is chosen as for the pointwise bands, whereas c_α is chosen as the $1 - \alpha$ quantile of $K_{\Phi,a}$.

Quantile-quantile plots are plots of $\widehat{x} = \widehat{F}^{-1}(u)$ against $x = F^{-1}(u)$ on a $\mathbf{R} \times \mathbf{R}$ -

square. Test bands of the type $u \pm n^{-1/2}c_\alpha\sigma_u$ can be constructed. Two different bands are worth noting.

Pointwise test bands can be established directly from Theorem 2.7 and (3.4). Here c_α is the $1 - \alpha/2$ quantile of the standard normal distribution and σ_u^2 equals $\text{Var}[\mathbb{X}_F\{F(x)\}]/\{f(x)\}^2$ at the point x . For the standard normal case, $F = \Phi$, the variance is reported in Theorem 2.9.

Kolmogorov-Smirnov bands can be constructed by $\sigma_u = 1$ and choosing c_α from the $1 - \alpha$ quantile of R_F .

3.4 Test for normality based on cumulants

Assuming an intercept is included in the model, the empirical cumulants can be computed from the sample moments of the standardised residuals, so

$$\hat{\kappa}_3 = \hat{\mu}_3, \quad \hat{\kappa}_4 = \hat{\mu}_4 - 3.$$

The following result follows directly from Corollary 2.10.

Corollary 3.1 *Suppose the model (2.3) and the Assumptions 2.1, 2.2, 2.3, 2.5 are satisfied, and that the innovations ε_t are $\mathbf{N}[0, 1]$ -distributed so $F(u) = \Phi(u)$. Then*

$$\left(\hat{\kappa}_3\sqrt{T/6}, \hat{\kappa}_4\sqrt{T/24}\right) \xrightarrow{D} \mathbf{N}_2[0, I_2].$$

Note that Assumption 2.2 is trivially satisfied, since the cumulants are based on the central moments, regardless of whether an intercept is included in the model. The Theorem presented here requires that the innovations ε_t are independent as stated in Assumption 2.4. This assumption is needed to be able to work with the empirical process, but could be relaxed with a proof based on direct expansions of the cumulants.

For models with independent observations, such a result was first established by Thiele (1889), see Lauritzen (2002) for a translation, and later by Pearson (1902), who argued that $(T/6)^{1/2}\hat{\kappa}_3$ and $(T/24)^{1/2}\hat{\kappa}_4$ are asymptotically standard normal. A large literature has investigated the joint behaviour in finite samples of independent observations. Jarque and Bera (1987) have given these tests a likelihood based motivation. Recently Kilian and Demiroglu (2002) have proved that these results hold for a class of cointegrated vector autoregressions with known cointegration rank. The proofs of these results are based on direct expansions of the sample moments of $\hat{\sigma}\hat{\varepsilon}_t$ rather than appealing to the empirical process of $\hat{\varepsilon}_t$. While the argument based on the empirical process requires independent innovations, an argument based on expansions of the sample moments could be made for innovations arising from a martingale difference sequence with constant conditional moments up to a certain order.

For cointegration applications the result of Kilian and Demiroglu (2002) can be combined with the present result in an interesting way. In a first step an unrestricted vector autoregression is fitted to the data and the hypothesis of normality is tested using the new result, where no knowledge of the autoregressive parameters is required. Secondly, a cointegration analysis is performed using the

likelihood procedure of Johansen (1995). Once the cointegration rank has been determined and imposed, the hypothesis of normality can then be tested once again for the restricted model using the results of Kilian and Demiroglu (2002).

3.5 Empirical illustration

To illustrate the methods a daily time series of average prices for whiting in the period Dec 1991 to May 1992 at the Fulton Fish Market was considered. This series was a part of a data set collected by Graddy (1995) with the purpose of estimating a demand function. The data are shown in Figure 1(a). A first order autoregression with an intercept was fitted giving the scaled residuals shown in panel (b). Before determining the location of the characteristic root, let alone estimating a demand function, the model assumptions need to be checked. This can be done without knowledge of the characteristic roots. The lag length can be determined using likelihood based tests or information criteria, see Nielsen (2006b) and normality can be checked using the methods developed in this paper.

Panels (c) and (d) show P-P and Q-Q plots as described above. For the P-P plot the 95% critical value for the two-sided Kolmogorov-Smirnov test is 0.903 according to Table 1. For the Q-Q plot the 95% critical value for the Kolmogorov-Smirnov-type statistic over the interval $5\% < p < 95\%$ is found to be 3.79 using the response surface for $R_{\Phi,0.45}$ in Table 2. Similarly the 95% critical value for the Anderson-Darling-type statistics for P-P and Q-Q plots are based on a critical value of 3.41 found from the response surface for $K_{\Phi,0.45}$. It is interesting to note that the Kolmogorov-Smirnov test is associated with P-P plots with bands that are parallel to the 45° line, thus having very little power against deviations in the tails. There does not appear to be any theoretical guidance towards choosing the cut-off points for the Anderson-Darling statistics and the Kolmogorov-Smirnov statistics for the Q-Q plot.

4 Proofs

4.1 Proof of main theorem

Theorem 2.7 is now proved. For the asymptotic analysis, it is convenient to decompose the empirical process. To facilitate this, the set $(\widehat{\varepsilon}_t \leq x)$ is rewritten in three steps: First, both sides of the inequality are scaled by $\widehat{\sigma}/\sigma$ to bring the residuals to the population scale; secondly, $\varepsilon_t - \widehat{\sigma}\widehat{\varepsilon}_t/\sigma$ is added to both sides; and thirdly, x is added and subtracted on the right. This gives

$$(\widehat{\varepsilon}_t \leq x) = (\varepsilon_t \leq x + \widehat{z}_t), \quad (4.1)$$

where

$$\widehat{z}_t = \widehat{a}x + \widehat{b}_t, \quad \text{where} \quad \widehat{a} = \frac{\widehat{\sigma}}{\sigma} - 1, \quad \widehat{b}_t = \varepsilon_t - \frac{\widehat{\sigma}}{\sigma}\widehat{\varepsilon}_t. \quad (4.2)$$

The empirical distribution of the residuals can then be decomposed as

$$\widehat{\mathbb{F}}(u) = \widehat{\mathbb{U}}(u) + \widehat{\mathbb{V}}(u) + \widehat{\mathbb{W}}(u), \quad (4.3)$$

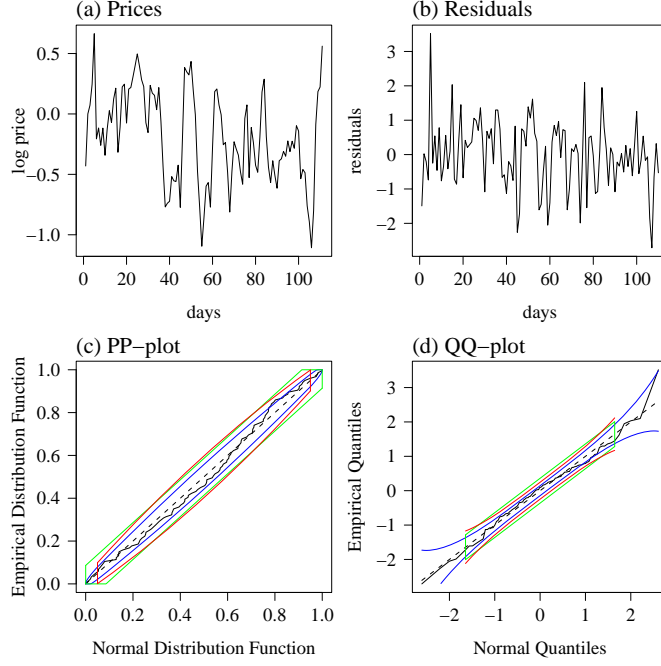


Figure 1: In panel (c) the point wise confidence band is shown with blue, the Kolmogorov-Smirnov band with green, and the Anderson-Darling band for $5\% < p < 95\%$ with red. In panel (d) the point wise confidence band is shown with blue, the Kolmogorov-Smirnov-type with green, and the Anderson-Darling-type band with red, both for $5\% < p < 95\%$.

where

$$\begin{aligned}\widehat{U}(u) &= \frac{1}{\sqrt{T}} \sum_{t=1}^T [1_{\{\varepsilon_t \leq F^{-1}(u)\}} - u], \\ \widehat{V}(u) &= f\{F^{-1}(u)\} \frac{1}{\sqrt{T}} \sum_{t=1}^T \widehat{z}_t, \\ \widehat{W}(u) &= \frac{1}{\sqrt{T}} \sum_{t=1}^T [1_{\{\widehat{\varepsilon}_t \leq F^{-1}(u)\}} - 1_{\{\varepsilon_t \leq F^{-1}(u)\}} - f\{F^{-1}(u)\} \widehat{z}_t].\end{aligned}$$

The components \widehat{U} and \widehat{V} are the leading terms. They are analysed in Theorem 4.2 below. For the analysis of \widehat{V} it is convenient to decompose $\widehat{V} = \widehat{V}_1 + \widehat{V}_2$, using the definition of \widehat{z}_t in (4.1). First, let

$$\widehat{V}_1(u) = f\{F^{-1}(u)\} \frac{1}{\sqrt{T}} \sum_{t=1}^T \left(\varepsilon_t - \frac{\widehat{\sigma}}{\sigma} \widehat{\varepsilon}_t \right) = f\{F^{-1}(u)\} \frac{1}{\sqrt{T}} \sum_{t=1}^T \varepsilon_t, \quad (4.4)$$

since $\sum_{t=1}^T \widehat{\varepsilon}_t = 0$ when an intercept is included in the model, as stipulated in Assumption 2.2, whereas, letting $F(x) = u$,

$$\widehat{V}_2(u) = F^{-1}(u) f\{F^{-1}(u)\} \sqrt{T} \left(\frac{\widehat{\sigma}}{\sigma} - 1 \right). \quad (4.5)$$

The third component, \widehat{W} , vanishes. For the analysis of \widehat{W} it is convenient to decompose $\widehat{W} = \widehat{W}_1 + \widehat{W}_2 + \widehat{W}_3$, letting $F(x) = u$,

$$\begin{aligned} \widehat{W}_1(u) &= \frac{1}{\sqrt{T}} \sum_{t=1}^T \{1_{(\varepsilon_t \leq x + \widehat{a}x)} - 1_{(\varepsilon_t \leq x)} - f(x) \widehat{a}x\}, \\ \widehat{W}_2(u) &= \frac{1}{\sqrt{T}} \sum_{t=1}^T \{F(x + \widehat{a}x + \widehat{b}_t) - F(x + \widehat{a}x) - f(x) \widehat{b}_t\} \\ \widehat{W}_3(u) &= \frac{1}{\sqrt{T}} \sum_{t=1}^T \{1_{(\varepsilon_t \leq x + \widehat{a}x + \widehat{b}_t)} - 1_{(\varepsilon_t \leq x + \widehat{a}x)} - F(x + \widehat{a}x + \widehat{b}_t) + F(x + \widehat{a}x)\}. \end{aligned}$$

The first term deals with the estimation of the scale parameter σ , while the other two deal with the estimation of expectation parameters. The first and the third term are referred to as uniform asymptotic linearity properties by Koul (2002), whereas the second is a Taylor expansion. It is shown in Theorems 4.3, 4.5, 4.6 that these terms disappear. The proof of Theorem 2.7 can now be summarised as follows.

Proof of Theorem 2.7. Decompose $\widehat{F} = \widehat{U} + \widehat{V}_1 + \widehat{V}_2 + \widehat{W}_1 + \widehat{W}_2 + \widehat{W}_3$. Theorem 4.2 shows the convergence of the leading terms $\widehat{U}, \widehat{V}_1, \widehat{V}_2$. Theorems 4.3, 4.5, 4.6 show that $\widehat{W}_1, \widehat{W}_2, \widehat{W}_3$ vanish. ■

Before looking at the asymptotic behaviour of \widehat{U} and \widehat{V} it is convenient to consider the estimators for the covariance parameters.

Lemma 4.1 *It holds that*

$$\frac{\widehat{\sigma}}{\sigma} - 1 = \frac{1}{2} \left(\frac{\widehat{\sigma}^2}{\sigma^2} - 1 \right) + O \left\{ \left(\frac{\widehat{\sigma}^2}{\sigma^2} - 1 \right)^2 \right\}.$$

Suppose model (2.3) and Assumptions 2.1, 2.3 are satisfied. Then

$$\sqrt{T} \left(\frac{\widehat{\sigma}^2}{\sigma^2} - 1 \right) = \frac{1}{\sqrt{T}} \sum_{t=1}^T (\varepsilon_t^2 - 1) + o_P(1) = O_P(1), \quad (4.6)$$

as well as $\sqrt{T}(\widehat{\rho} - \rho) = O_P(1)$.

Proof of Lemma 4.1. The first result follows from the Taylor expansion

$$\frac{\widehat{\sigma}}{\sigma} - 1 = \sqrt{1 + \frac{\widehat{\sigma}^2 - \sigma^2}{\sigma^2}} - 1 = \frac{\widehat{\sigma}^2 - \sigma^2}{2\sigma^2} + O_P \left\{ \left(\frac{\widehat{\sigma}^2 - \sigma^2}{\sigma^2} \right)^2 \right\}.$$

Next, Nielsen (2005, Theorem 2.6) shows that under Assumptions 2.1, 2.3, noting that $\lambda > 4$ in (2.6), it holds

$$\widehat{\Omega} \stackrel{a.s.}{=} \frac{1}{T} \sum_{t=1}^T \xi_t \xi_t' + o(T^{-1/2}).$$

The Central Limit Theorem for martingale differences by Brown and Eagleson (1971) is applicable when (2.6), (2.7) are satisfied, and implies that $\sqrt{T}(\widehat{\Omega} - \Omega)$ is asymptotically normal. Using the functional δ -method it is seen that this property is shared by $\widehat{\sigma}$ and $\widehat{\rho}$. ■

Theorem 4.2 *Suppose model (2.3) and Assumption 2.1, 2.2, 2.3, 2.4 are satisfied. Then, letting \odot denote the element wise (Hadamard) product,*

$$\begin{Bmatrix} \widehat{\mathbb{U}}(u) \\ \widehat{\mathbb{V}}_1(u) \\ \widehat{\mathbb{V}}_2(u) \end{Bmatrix} = \begin{bmatrix} 1 \\ \mathbf{f}\{\mathbb{F}^{-1}(u)\} \\ \frac{1}{2}\mathbb{F}^{-1}(u) \mathbf{f}\{\mathbb{F}^{-1}(u)\} \end{bmatrix} \odot \int_0^1 \begin{bmatrix} 1_{(s \leq u)} \\ \mathbb{F}^{-1}(s) \\ \{\mathbb{F}^{-1}(s)\}^2 \end{bmatrix} d\widehat{\mathbb{U}}(s) + \begin{Bmatrix} 0 \\ 0 \\ \text{OP}(1) \end{Bmatrix},$$

which converges in distribution to $\mathbb{X}_{\mathbb{F}}(u)$ on $\mathbb{D}[0, 1]$.

Proof of Lemma 4.2. First, the stochastic integral representation in terms of $\widehat{\mathbb{U}}$ is established. For the first component this follows from the discussion in §2.1. For the other two terms, denote the empirical distribution of the innovations by $\mathbb{F}_T(x) = T^{-1} \sum_{t=1}^T 1_{(\varepsilon_t \leq x)}$. As in §2.3 the sample and population moments of ε_t satisfy

$$\begin{aligned} \frac{1}{\sqrt{T}} \sum_{t=1}^T (\varepsilon_t^m - \mathbb{E}\varepsilon_t^m) &= \sqrt{T} \int_{\mathbf{R}} x^m d\{\mathbb{F}_T(x) - \mathbb{F}(x)\} \\ &= \int_{\mathbf{R}} x^m d\widehat{\mathbb{U}}\{\mathbb{F}(x)\} = \int_0^1 \{\mathbb{F}^{-1}(u)\}^m d\widehat{\mathbb{U}}(u), \quad (4.7) \end{aligned}$$

substituting $u = \mathbb{F}(x)$. Thus the expression for $\widehat{\mathbb{V}}_1$, follows directly (4.4), which requires Assumption 2.2, whereas the expression for $\widehat{\mathbb{V}}_2$ follows from (4.5) and Lemma 4.1, requiring the Assumptions 2.1, 2.3.

Billingsley (1968, §13) shows that a continuous $\widehat{\mathbb{U}}$ converges to \mathbb{U} on $\mathbb{C}[0, 1]$. This requires Assumption 2.4. Since the limit is continuous, $\widehat{\mathbb{U}}$ converges to \mathbb{U} on $\mathbb{D}[0, 1]$. Further, due to the existence of fourth moments, see Assumption 2.3, the integrand in (4.7) is square integrable. Following Shorack and Wellner (1986, p.94) the integrator $\widehat{\mathbb{U}}$ can be replaced with the Brownian bridge \mathbb{U} in the limit. This gives the process $\mathbb{X}_{\mathbb{F}}$ in (2.9). ■

It is now argued that the component $\widehat{\mathbb{W}}_1$ vanishes. This result follows from the work of Koul (2002).

Theorem 4.3 *Suppose model (2.3) and Assumptions 2.1, 2.3, 2.6 are satisfied. Then*

$$\sup_{0 \leq u \leq 1} \left| \widehat{\mathbb{W}}_1(u) \right| = o_{\mathbf{P}}(1).$$

Proof. From Lemma 4.1 it follows that \widehat{a} is $O_{\mathbf{P}}(T^{-1/2})$. As pointed out by Rao and Sethuraman (1975) and Loynes (1980) a bound $b > 0$ can be found so $|\widehat{a}| < b$ with probability close to one. Thus it suffices to show

$$\sup_{\substack{x \in \mathbf{R} \\ |s| < b}} \frac{1}{\sqrt{T}} \sum_{t=1}^T \left[1_{\{\varepsilon_t \leq x(1+T^{-1/2}s)\}} - 1_{\{\varepsilon_t \leq x\}} - f(x) T^{-1/2} s x \right].$$

This result follows from Corollary 2.3.2 of Koul (2002, p.59). A set of assumptions have to be checked. First, note that, in the notation of Koul (2002), $n = T$, $d_{ni} = T^{-1/2}$, $\mathbf{c}_{ni} = 0$, $X_{ni} = \varepsilon_t$, $H(x) = F_{ni} = F(x)$, $f_{ni}(x) = f(x)$. Since d_{ni} is uniform in t , the conditions to d_{ni} in **N1** and **N2** of Koul (2002, p.16) are trivially satisfied. Since $\mathbf{c}_{ni} = 0$ then the conditions to \mathbf{c}_{ni} in (2.3.6) and (2.3.7) of Koul (2002, p.52) are trivially satisfied. The conditions **F1**, **F2**, **F3** to f of Koul (2002, p.59) are satisfied by Assumption 2.6 as follows: **F1** requires uniform continuity of f , which is satisfied since f is differentiable and $\sup_{x \in \mathbf{R}} |f'(x)| < \infty$; **F2** requires f to be positive; **F3** requires $\sup_{x \in \mathbf{R}} |xf(x)| < \infty$. ■

Before looking at the components $\widehat{\mathbb{W}}_2$ and $\widehat{\mathbb{W}}_3$ it is useful to discuss the properties of the process $\widehat{z}_t = \widehat{a}x + \widehat{b}_t$ defined in (4.2).

Lemma 4.4 *Suppose model (2.3) and Assumptions 2.1, 2.3 are satisfied. Then, if $g(T) \rightarrow \infty$ as $T \rightarrow \infty$,*

(i) $\sigma \widehat{b}_t = (0, \widehat{\rho} - \rho) \xi_t + (1, -\widehat{\rho}) (\widehat{\theta} - \theta) S_{t-1}$,

(ii) $T^{-1/2} \sum_{t=1}^T \widehat{z}_t^2 = (1 + x^2) o_{\mathbf{P}}(1)$,

(iii) \widehat{b}_t can be written as $\widehat{b}_t = -\boldsymbol{\alpha}_T \mathbf{z}_{Tt}$ where

(a) \mathbf{z}_{Tt} is \mathcal{G}_{t-1} -measurable as defined in Assumption 2.5,

(b) $\boldsymbol{\alpha}_T = O_{\mathbf{P}}(1)$,

(c) $\sum_{t=1}^{T-g(T)} \mathbf{z}_{Tt} \mathbf{z}'_{Tt} = O_{\mathbf{P}}(1)$ for any function g so $g(T)/\log T \rightarrow \infty$.

Proof of Lemma 4.4. (i): rewrite \widehat{b}_t in terms of the parameters of the vector autoregression (2.3). By definition $\sigma \widehat{b}_t = \sigma \varepsilon_t - \widehat{\sigma} \widehat{\varepsilon}_t$, where $\sigma \varepsilon_t = (1, -\rho) \xi_t$ and $\widehat{\sigma} \widehat{\varepsilon}_t = (1, -\widehat{\rho}) \widehat{\xi}_t$, and ρ is either 0 or $\Omega_{yz} \Omega_{zz}^{-1}$. Adding and subtracting $(1, -\widehat{\rho}) \xi_t$ gives

$$\sigma b_t = \sigma \varepsilon_t - \widehat{\sigma} \widehat{\varepsilon}_t = (1, -\rho) \xi_t - (1, -\widehat{\rho}) \widehat{\xi}_t = (0, \widehat{\rho} - \rho) \xi_t - (1, -\widehat{\rho}) (\widehat{\xi}_t - \xi_t).$$

Writing the vector autoregression (2.3) in companion form

$$X_t = \theta S_{t-1} + \xi_t \quad \text{where} \quad \theta = (A_1, \dots, A_k, \mu), \quad S_{t-1} = (X'_{t-1}, \dots, X'_{t-k}, D'_{t-1})'$$

and in particular $\widehat{\xi}_t - \xi_t = -(\widehat{\theta} - \theta) S_{t-1}$ leads to the desired expression.

(ii): The least squares estimator for θ satisfies $\widehat{\theta} - \theta = \mathbf{M}_{\xi S} \mathbf{M}_{SS}^{-1}$ where

$$\mathbf{M}_{\xi\xi} = \sum_{t=1}^T \xi_t \xi_t' \quad \mathbf{M}_{\xi S} = \sum_{t=1}^T \xi_t S_{t-1}', \quad \mathbf{M}_{SS} = \sum_{t=1}^T S_{t-1} S_{t-1}'.$$

Using the inequality $(x + y)^2 \leq 2(x^2 + y^2)$ both to \widehat{z}_t^2 and to $(\widehat{\sigma b}_t)^2$ shows

$$\begin{aligned} \frac{1}{4} \sum_{t=1}^T (\widehat{\sigma z}_t)^2 &\leq \frac{T}{2} (\widehat{\sigma a x})^2 + \frac{1}{2} \sum_{t=1}^T (\widehat{\sigma b}_t)^2 \\ &\leq T (\widehat{\sigma a x})^2 + (0, \widehat{\rho} - \rho) \mathbf{M}_{\xi\xi} (0, \widehat{\rho} - \rho)' + (1, -\widehat{\rho}) \mathbf{M}_{\xi S} \mathbf{M}_{SS}^{-1} \mathbf{M}_{S\xi} (1, -\widehat{\rho})'. \end{aligned}$$

From Lemma 4.1, it follows that \widehat{a} and $\widehat{\rho} - \rho$ are both $\mathcal{O}_{\mathbb{P}}(T^{-1/2})$. Further, Nielsen (2005, Theorem 2.4, 6.1) shows that $\mathbf{M}_{\xi S} \mathbf{M}_{SS}^{-1} \mathbf{M}_{S\xi} = \mathcal{O}_{\mathbb{P}}(T^{1/2})$ and $\mathbf{M}_{\xi\xi} = \mathcal{O}_{\mathbb{P}}(1)$. Both arguments use Assumptions 2.1, 2.3. Normalising by $T^{-1/2}$ gives the desired result.

(iii): From (i) it holds that $\widehat{\sigma b}_t = \sigma \boldsymbol{\alpha}_T \mathbf{z}_{Tt}$, since, for instance in the presence of ρ ,

$$\boldsymbol{\alpha}_T = - \left\{ (\widehat{\rho} - \rho), (1, -\widehat{\rho}) (\widehat{\theta} - \theta) \right\} N_T, \quad \mathbf{z}_{Tt} = N_T^{-1} \left\{ \begin{array}{c} (0, I_{p-1}) \xi_t \\ S_{t-1} \end{array} \right\},$$

for some normalisation matrix N_T . By construction \mathbf{z}_{Tt} is \mathcal{G}_{t-1} -measurable, showing (a). Note, that (ii) implies that $\sum_{t=1}^T \boldsymbol{\alpha}_T \mathbf{z}_{Tt} \mathbf{z}_{Tt}' \boldsymbol{\alpha}_T' = \mathcal{O}_{\mathbb{P}}(T^{1/2})$, which is not quite sufficient for (b), (c). Thus, use that a matrix M exists so

$$M S_t = \begin{pmatrix} U_t \\ Q_t \\ W_t \end{pmatrix} = \begin{pmatrix} \mathbf{U} & 0 & 0 \\ 0 & \mathbf{Q} & 0 \\ 0 & 0 & \mathbf{W} \end{pmatrix} \begin{pmatrix} U_{t-1} \\ Q_{t-1} \\ W_{t-1} \end{pmatrix} = \begin{pmatrix} e_{U,t} \\ e_{Q,t} \\ e_{W,t} \end{pmatrix},$$

where the absolute values of the eigenvalues of \mathbf{U} , \mathbf{Q} , \mathbf{W} are less than one, equal to one, and greater than one, respectively, see Nielsen (2005, §3). The deterministic components are therefore included in the Q_t process. Using Theorems 2.4, 6.2 and Table 1 of Nielsen (2005) it follows that

$$\begin{pmatrix} \mathbf{M}_{\xi\xi} & \mathbf{M}_{\xi S} \\ \mathbf{M}_{S\xi} & \mathbf{M}_{SS} \end{pmatrix} = \text{diag}(\mathbf{M}_{\xi\xi}, \mathbf{M}_{UU}, \mathbf{M}_{QQ}, \mathbf{M}_{WW}) \{1 + \mathcal{O}_{\mathbb{P}}(1)\}. \quad (4.8)$$

Therefore, write

$$\widehat{\sigma b}_t = -\sigma (\boldsymbol{\alpha}_{\xi, T} \mathbf{z}_{\xi, Tt}, \boldsymbol{\alpha}_{U, T} \mathbf{z}_{U, Tt}, \boldsymbol{\alpha}_{Q, T} \mathbf{z}_{Q, Tt}, \boldsymbol{\alpha}_{W, T} \mathbf{z}_{W, Tt}) \{1 + \mathcal{O}_{\mathbb{P}}(1)\},$$

where, for $t \leq T - g(T)$ and some normalisation matrix $N_{Q, T}$,

$$\begin{aligned} \sigma \boldsymbol{\alpha}_{\xi, t} &= -(\widehat{\rho} - \rho) T^{1/2}, & \mathbf{z}_{\xi, Tt} &= T^{-1/2} (0, I_{p-1}) \xi_t, \\ \sigma \boldsymbol{\alpha}_{U, t} &= -(1, -\widehat{\rho}) \mathbf{M}_{\xi U} \mathbf{M}_{UU}^{-1} T^{1/2}, & \mathbf{z}_{U, Tt} &= T^{-1/2} U_{t-1}, \\ \sigma \boldsymbol{\alpha}_{Q, t} &= -(1, -\widehat{\rho}) \mathbf{M}_{\xi Q} \mathbf{M}_{QQ}^{-1} N_{Q, T}, & \mathbf{z}_{Q, Tt} &= N_{Q, T}^{-1} Q_{t-1}, \\ \sigma \boldsymbol{\alpha}_{W, t} &= -(1, -\widehat{\rho}) \mathbf{M}_{\xi W} \mathbf{M}_{WW}^{-1} \mathbf{W}^{T-g(T)}, & \mathbf{z}_{W, Tt} &= \mathbf{W}^{g(T)-T} W_{t-1}. \end{aligned}$$

For (b) it needs to be argued that each α -term is $O_{\mathbb{P}}(1)$. For (c) it suffices to show that each \mathbf{z} -term has sums of squares that are $O_{\mathbb{P}}(1)$, and that the cross-products are $o_{\mathbb{P}}(1)$. For each of the following arguments it suffices that Assumptions 2.1, 2.3 are satisfied.

The ξ terms. Lemma 4.1 shows that $\alpha_{\xi,T} = O_{\mathbb{P}}(1)$. Nielsen (2005, Theorem 6.2) shows that $\sum_{t=1}^T \mathbf{z}_{\xi,Tt} \mathbf{z}'_{\xi,Tt}$ is $O_{\mathbb{P}}(1)$.

The U term. The Central Limit Theorem for martingale differences by Brown and Eagleson (1971) implies that $\mathbf{M}_{\xi U} \mathbf{M}_{UU}^{-1/2} = O_{\mathbb{P}}(1)$. Nielsen (2005, Theorem 6.2) shows that $T^{-1} \mathbf{M}_{UU}$ has a positive definite limit, *a.s.* Thus $\alpha_{U,T}$, $\sum_{t=1}^T \mathbf{z}_{U,Tt} \mathbf{z}'_{U,Tt}$ are both $O_{\mathbb{P}}(1)$.

The Q term. Using standard unit root weak convergence arguments as in Chan and Wei (1988) and Chan (1989) it can be shown that a normalisation matrix $N_{Q,T}$ exists so $\mathbf{M}_{\xi Q} N_{Q,T}$ and $N_{Q,T}^{-1} \sum_{t=1}^T Q_{t-1} Q'_{t-1} (N_{Q,T}^{-1})'$ are $O_{\mathbb{P}}(1)$, and the latter is positive definite *a.s.* Thus $\alpha_{Q,T}$, $\sum_{t=1}^T \mathbf{z}_{Q,Tt} \mathbf{z}'_{Q,Tt}$ have the desired properties. Indeed Chan and Wei (1988), consider the univariate case, where $p = 1$ and $X_t = Y_t$, without deterministic terms, whereas Chan (1989) include deterministic terms for that case. The idea of Chan and Wei (1988) is first to show in their Theorem 2.2 that $T^{-1/2} \sum_{t=1}^{\text{int}(Tu)} \mathbf{\Lambda}_t \varepsilon_t$ converges to vector of independent standard Brownian motions, when defining

$$\mathbf{\Lambda}_t = \{1, (-1)^t, \sqrt{2} \sin(t\theta_1), \sqrt{2} \cos(t\theta_1), \dots, \sqrt{2} \sin(t\theta_l), \sqrt{2} \cos(t\theta_l)\},$$

for $\theta_k \in (0, \pi)$ so $\theta_k \neq \theta_j$ if $k \neq j$. This results is easily generalised to a multivariate result using the Cramér-Wold device, see Billingsley (1968). Next, existence of a normalisation matrix $N_{Q,T}$ and convergence of $T^{1/2} N_{Q,T}^{-1} Q_t$ and $N_{Q,T}^{-1} \sum_{t=1}^T Q_{t-1} Q'_{t-1} (N_{Q,T}^{-1})'$ then follows from the Continuous Mapping Theorem (see Chan and Wei, 1988, §3), which is therefore easily generalisable. For $\mathbf{M}_{\xi Q} N_{Q,T}$, convergence of $\mathbf{M}_{\xi Q} N_{Q,T}$ jointly with that of $T^{1/2} N_{Q,T}^{-1} Q_t$ is shown by a Skorokhod embedding result when \mathbf{Q} is univariate (Chan and Wei, 1988, Theorem 2.4(ii)). This is easy to generalise when \mathbf{Q} is multivariate with simple roots (Chan and Wei, 1988, Remark to Theorem 2.4). For general \mathbf{Q} the Continuous Mapping Theorem is used (Chan and Wei, 1988, Theorem 2.4(i)), which again is easily generalisable. Finally, the positive definiteness of $N_{Q,T}^{-1} \sum_{t=1}^T Q_{t-1} Q'_{t-1} (N_{Q,T}^{-1})'$ is shown (Chan and Wei, 1988, Lemma 3.1.1), which can be generalised using a Cramér-Wold-type argument.

The W term. Nielsen (2005, Theorem 2.4, Corollary 7.2) shows $\mathbf{M}_{\xi W} \mathbf{M}_{WW}^{-1/2} = o(T^{1/4})$, *a.s.* and that $\mathbf{W}^{-T} \mathbf{M}_{WW} (\mathbf{W}^{-T})'$ is convergent with positive definite limiting points. In particular $\sum_{t=1}^T \mathbf{z}_{W,Tt} \mathbf{z}'_{W,Tt} = \mathbf{W}^{g(T)-T} \sum_{t=1}^{T-g(T)} W_{t-1} W'_{t-1} (\mathbf{W}^{g(T)-T})'$ is convergent with positive definite limiting points, while $-\sigma \alpha_{W,T} = o(T^{1/4} \mathbf{W}^{-g(T)})$ *a.s.*, which again is $o(1)$ for any $g(T)$ so $g(T)/\log T \rightarrow \infty$.

Cross terms. These all vanish. This follows directly from (4.8). ■

Theorem 4.5 *Suppose model (2.3) and Assumptions 2.1, 2.3, 2.6 are satisfied. Then*

$$\sup_{0 \leq u \leq 1} \left| \widehat{\mathbb{W}}_2(u) \right| = o_{\mathbb{P}}(1).$$

Proof of Theorem 4.5. At first $\widehat{\mathbb{W}}_2(u)$ is written as an integral

$$\widehat{\mathbb{W}}_2(u) = \frac{1}{\sqrt{T}} \sum_{t=1}^T \{F(x + \widehat{z}_t) - F(x) - f(x) \widehat{z}_t\} = \frac{1}{\sqrt{T}} \sum_{t=1}^T \int_x^{x+\widehat{z}_t} \{f(y) - f(x)\} dy.$$

By the triangle inequality

$$\left| \widehat{\mathbb{W}}_2(u) \right| \leq \frac{1}{\sqrt{T}} \sum_{t=1}^T \int_x^{x+\widehat{z}_t} |f(y) - f(x)| dy.$$

The integrand can be bounded by its maximum, so

$$\left| \widehat{\mathbb{W}}_2(u) \right| \leq \frac{1}{\sqrt{T}} \sum_{t=1}^T |\widehat{z}_t| \max_{|h| \leq |\widehat{z}_t|} |f(x+h) - f(x)|.$$

The mean value theorem then implies a further bound

$$\left| \widehat{\mathbb{W}}_2(u) \right| \leq \frac{2}{\sqrt{T}} \sum_{t=1}^T |\widehat{z}_t|^2 \max_{|h| \leq |\widehat{z}_t|} |f'(x+h)|.$$

Taking the maximum over the entire real axis, and using Lemma 4.4(ii), which requires Assumptions 2.1, 2.3, gives

$$\left| \widehat{\mathbb{W}}_2(u) \right| \leq o_{\mathbb{P}}(1) \sup_{x \in \mathbf{R}} |(1+x^2) f'(x)|,$$

which follows from Assumption 2.6. ■

The asymptotic uniform linearity property, that $\widehat{\mathbb{W}}_3$ vanishes, is now proved. Two ideas of Lee and Wei (1999) are used. First, to deal with the issue that for explosive component $W_T W_T'$ and $\sum_{t=1}^T W_t W_t'$, the largest and the smallest components are treated separately. Lee and Wei do actually not consider explosive and non-explosive components jointly, but the joint evaluation turns out to not to pose any problems. Secondly, Theorem 2.2 of Lee and Wei gives an asymptotic uniform linearity property for triangular arrays, which can be used here.

Theorem 4.6 *Suppose model (2.3) and Assumptions 2.1, 2.3, 2.4, 2.5, 2.6 are satisfied. Then*

$$\sup_{0 \leq u \leq 1} \left| \widehat{\mathbb{W}}_3(u) \right| = o_{\mathbb{P}}(1).$$

Proof of Theorem 4.6. Some notation is needed. Let $u = F(x)$ and, recalling $\widehat{a}_t, \widehat{b}_t$ defined in (4.2), let $\widehat{x} = x(1 + \widehat{a}_t)$. Define

$$\mathbf{w}(t, \widehat{x}) = 1_{(\varepsilon_t \leq \widehat{x} + \widehat{b}_t)} - 1_{(\varepsilon_t \leq \widehat{x})} - F(\widehat{x} + \widehat{b}_t) + F(\widehat{x}).$$

Decompose $\widehat{\mathbb{W}}_3 = \widehat{\mathbb{W}}_{3,1} + \widehat{\mathbb{W}}_{3,2}$ where

$$\widehat{\mathbb{W}}_{3,1}(u) = \frac{1}{\sqrt{T}} \sum_{t=1}^{T-g(T)} \mathbf{w}(t, \widehat{x}), \quad \widehat{\mathbb{W}}_{3,2}(\widehat{x}) = \frac{1}{\sqrt{T}} \sum_{t=T-g(T)+1}^T \mathbf{w}(t, \widehat{x}),$$

for some function $g(T)$ chosen so $g(T)/\sqrt{T} \rightarrow 0$ and $g(T)/\log T \rightarrow \infty$.

Analysis of $\widehat{\mathbb{W}}_{3,2}$. It is immediately seen that $|\mathbf{w}(t, u)| \leq 2$, so

$$\sup_{0 \leq u \leq 1} \left| \widehat{\mathbb{W}}_{3,2}(u) \right| \leq \frac{1}{\sqrt{T}} \sum_{t=T-g(T)+1}^T 2 = \frac{2g(T)}{\sqrt{T}} \rightarrow 0.$$

Analysis of $\widehat{\mathbb{W}}_{3,1}$. First, note that taking supremum over $x \in \mathbf{R}$ and over $\widehat{x} \in \mathbf{R}$ gives the same supremum so

$$\sup_{0 \leq u \leq 1} \left| \widehat{\mathbb{W}}_{3,1}(u) \right| = \sup_{x \in \mathbf{R}} \left| \frac{1}{\sqrt{T}} \sum_{t=1}^{T-g(T)} \mathbf{w}(t, \widehat{x}) \right| = \sup_{x \in \mathbf{R}} \left| \frac{1}{\sqrt{T}} \sum_{t=1}^{T-g(T)} \mathbf{w}(t, x) \right|.$$

Secondly, this can be written as

$$\sup_{x \in \mathbf{R}} \left| \frac{1}{\sqrt{T}} \sum_{t=1}^T \left\{ \mathbf{1}_{(\varepsilon_t \leq x - \boldsymbol{\alpha}_T \mathbf{z}_{Tt})} - \mathbf{1}_{(\varepsilon_t \leq x)} - \mathbf{F}(x - \boldsymbol{\alpha}_T \mathbf{z}_{Tt}) + \mathbf{F}(x) \right\} \right|,$$

where $-\boldsymbol{\alpha}_T \mathbf{z}_{Tt}$ equals \widehat{b}_t for $t \leq T-g(T)$ and zero otherwise. According to Lemma 4.4(iii) the term \mathbf{z}_{Tt} can be chosen as a triangular array of variables measurable with respect to the filtration \mathcal{G}_{t-1} introduced in Assumption 2.5.

Thirdly, the desired result now follows if the conditions of Lee and Wei (1999, Corollary 2.1) can be established. First, ε_t are independent and identically distributed with distribution function \mathbf{F} according to Assumption 2.4, and independent of \mathcal{G}_{t-1} by Assumption 2.5. The vectors $\boldsymbol{\alpha}'_T$ and \mathbf{z}_{Tt} have a dimension not depending on T where \mathbf{z}_{Tt} is \mathcal{G}_{t-1} -measurable. Since the ε_t s have the same marginal distribution function \mathbf{F} with uniformly bounded second derivatives by Assumption 2.6 then Lee and Wei's condition (2.11) is trivially satisfied. Finally $\boldsymbol{\alpha}_T$ and $\sum_{t=1}^T \mathbf{z}_{Tt} \mathbf{z}'_{Tt}$ are both $\text{O}_{\mathbf{P}}(1)$ according to Lemma 4.4(iii), requiring Assumption 2.1, 2.3 ■

4.2 Empirical moments: Proof of Theorem 2.8

The following steps are taken. First, it is argued that the sample moments converge in distribution. Secondly, the identities (2.12), (2.13) are proved. Thirdly, the asymptotic covariance matrix is derived.

First, by using Lemma 4.1 along with (4.7), (2.11) then

$$\begin{aligned} \sqrt{T} \left(\frac{\widehat{\sigma}^2}{\sigma^2} - 1 \right) &= \int_0^1 \{ \mathbf{F}^{-1}(u) \}^2 d\widehat{\mathbb{U}}(u), \\ \sqrt{T} (\widehat{\mu}_m - \mu_m) &= \int_0^1 \{ \mathbf{F}^{-1}(u) \}^m d\widehat{\mathbb{F}}(u). \end{aligned}$$

Since $\widehat{\mathbb{U}}$ is one of the components of $\widehat{\mathbb{F}}$, see (4.3), the joint convergence follows from Theorem 2.7, see Shorack and Wellner (1986, p.92). The second set of expressions for the limiting distribution in (2.12), (2.13) then follows. The first set of expressions will be derived below. The asymptotic, mean-zero, normality follows directly from the remarks about stochastic integrals with respect to \mathbb{U} in §2.2.

Secondly, the identity (2.12) is a straight forward definition. For the identity (2.13) some more work is required. The process $\mathbb{X}_{\mathbb{F}}$ is expressed in terms of \mathbb{U} in (2.9) showing that

$$\int_0^1 \{\mathbb{F}^{-1}(u)\}^m d\mathbb{X}_{\mathbb{F}}(u) = J_m + I_1 J_1 + I_2 J_2 \quad (4.9)$$

where the uniform process \mathbb{U} enters the J_m terms:

$$J_m = \int_0^1 \{\mathbb{F}^{-1}(u)\}^m d\mathbb{U}(u),$$

whereas the term I_1 and I_2 are standard integrals not involving \mathbb{U} , and which will be shown to equal:

$$\begin{aligned} I_1 &= \int_0^1 \{\mathbb{F}^{-1}(u)\}^m d[\mathfrak{f}\{\mathbb{F}^{-1}(u)\}] = -m\mu_{m-1}, \\ I_2 &= \int_0^1 \{\mathbb{F}^{-1}(u)\}^m d\left[\frac{1}{2}\mathbb{F}^{-1}(u)\mathfrak{f}\{\mathbb{F}^{-1}(u)\}\right] = -\frac{m}{2}\mu_m. \end{aligned}$$

The integral I_1 : The integrator is rewritten using the chain rule and implicit differentiation as

$$d[\mathfrak{f}\{\mathbb{F}^{-1}(u)\}] = \frac{\mathfrak{f}'\{\mathbb{F}^{-1}(u)\}}{\mathfrak{f}\{\mathbb{F}^{-1}(u)\}} du.$$

Inserting this in I_1 gives

$$I_1 = \int_0^1 \{\mathbb{F}^{-1}(u)\}^m \frac{\mathfrak{f}'\{\mathbb{F}^{-1}(u)\}}{\mathfrak{f}\{\mathbb{F}^{-1}(u)\}} du.$$

Substituting $u = \mathbb{F}(x)$ so $du = \mathfrak{f}(x) dx$ gives

$$I_1 = \int_{\mathbf{R}} x^m \frac{\mathfrak{f}'(x)}{\mathfrak{f}(x)} \mathfrak{f}(x) dx = \int_{\mathbf{R}} x^m \mathfrak{f}'(x) dx.$$

Finally, by partial integration

$$I_1 = -m \int_{\mathbf{R}} x^{m-1} \mathfrak{f}(x) dx = -m\mu_{m-1}.$$

The integral I_2 : In the same way

$$d\left[\frac{1}{2}\mathbb{F}^{-1}(u)\mathfrak{f}\{\mathbb{F}^{-1}(u)\}\right] = \frac{1}{2}\left[1 + \mathbb{F}^{-1}(u)\frac{\mathfrak{f}'\{\mathbb{F}^{-1}(u)\}}{\mathfrak{f}\{\mathbb{F}^{-1}(u)\}}\right] du.$$

Inserting this in I_2 , substituting $u = F(x)$ and using partial integration

$$I_2 = \frac{1}{2} \int_{\mathbf{R}} x^m \left\{ 1 + x \frac{f'(x)}{f(x)} \right\} f(x) dx = \frac{1}{2} \{ \mu_m - (m+1) \mu_m \} = -\frac{m}{2} \mu_m.$$

Finally the covariance matrix: The covariance follow from the formula (2.8). Thus, the expression for σ_{22} follows from

$$\int_0^1 h_2^2(u) du = \int_0^1 \{F^{-1}(u)\}^4 du = \mu_4, \quad \int_0^1 h_2(u) du = \mu_2 = 1.$$

The other expressions follow in a similar way. First,

$$\int_0^1 h_m(u) du = \mu_m - m\mu_{m-1}\mu_1 - \frac{m}{2}\mu_m\mu_2 = \mu_m \left(1 - \frac{m}{2} \right),$$

using $\mu_1 = 0$ and $\mu_2 = 1$, while

$$\int_0^1 h_m(u) h_2(u) du = \mu_{m+2} - m\mu_{m-1}\mu_3 - \frac{m}{2}\mu_m\mu_4$$

cannot be reduced further. Combining this with $\int_0^1 h_2(u)du$ and $\int_0^1 h_m(u)du$ leads to σ_{m2} . Finally, σ_{mn} can be derived by combining $\int_0^1 h_m(u)du$ with

$$\begin{aligned} \int_0^1 h_m(u) h_n(u) du &= \mu_{m+n} - n\mu_{m+1}\mu_{n-1} - \frac{n}{2}\mu_{m+2}\mu_n \\ &\quad - m\mu_{m-1}\mu_{n+1} + mn\mu_{m-1}\mu_{n-1}\mu_2 + \frac{mn}{2}\mu_{m-1}\mu_n\mu_3 \\ &\quad - \frac{m}{2}\mu_m\mu_{n+2} + \frac{mn}{2}\mu_m\mu_{n-1}\mu_3 + \frac{mn}{4}\mu_m\mu_n\mu_4. \end{aligned}$$

4.3 The Gaussian case: Proof of Theorem 2.9

It is first established, for $\Phi(x) = u$ that

$$\begin{aligned} \int_0^1 1_{(s \leq u)} \Phi^{-1}(s) ds &= -\varphi(x), \\ \int_0^1 1_{(s \leq u)} \{ \Phi^{-1}(s) \}^2 ds &= -x\varphi(x) + u. \end{aligned}$$

To see this, note that $\varphi'(x) = -x\varphi(x)$ and $\varphi''(x) = (x^2 - 1)\varphi(x)$, and therefore by substitution $\Phi(t) = s$ it holds

$$\int_0^1 1_{(s \leq u)} \Phi^{-1}(s) ds = \int_{-\infty}^x t\varphi(t) dt = - \int_{-\infty}^x \varphi'(t) dt = -\varphi(x),$$

and likewise

$$\begin{aligned} \int_0^1 1_{(s \leq u)} \{ \Phi^{-1}(s) \}^2 ds &= \int_{-\infty}^x \{ (t^2 - 1) + 1 \} \varphi(t) dt \\ &= \int_{-\infty}^x \{ \varphi''(t) + \varphi(t) \} dt = \varphi'(x) + \Phi(x). \end{aligned}$$

The integrals involved in the expression (2.8) for the covariance of integrals with respect to the Brownian bridge can now be considered. Recalling the integral expressions for the Gaussian moments in (2.14) it holds with

$$g_u(s) = \left[1_{(s \leq u)}, \Phi^{-1}(s), \{\Phi^{-1}(s)\}^2 \right]',$$

that

$$\varkappa_u = \int_0^1 g_u(s) ds = (u, 0, 1)',$$

and for $u = \Phi(x)$ and $v = \Phi(y)$ so $u \leq v$

$$\Sigma_{u,v} = \int_0^1 g_u(s) \{g_v(s)\}' ds = \begin{pmatrix} u & -\varphi(x) & u - x\varphi(x) \\ -\varphi(y) & 1 & 0 \\ u - y\varphi(y) & 0 & 3 \end{pmatrix}.$$

The desired expression is then computed as

$$\{1, \varphi(x), x\varphi(x)/2\} (\Sigma_{u,v} - \varkappa_u \varkappa_v') \{1, \varphi(y), y\varphi(y)/2\}'.$$

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6 References

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