# Identification-Robust Estimation and Testing of the Zero-Beta CAPM

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We propose exact simulation-based procedures for: (i) testing mean-variance efficiency when the zero-beta rate is unknown; (ii) building confidence intervals for the zero-beta rate. On observing that this parameter may be weakly identified, we propose likelihood-ratio-type tests as well as Fieller-type procedures based on a Hotelling-HAC statistic, which are robust to weak identification and allow for non-Gaussian distributions including parametric GARCH structures. The Fieller-Hotelling-HAC procedure also accounts (asymptotically) for general forms of heteroskedasticity and autocorrelation. We propose confidence sets for the zero-beta rate based on "inverting" exact tests for this parameter; for both procedures proposed, these sets can be interpreted as multivariate extensions of the classic Fieller method for inference on ratios. The exact distribution of likelihood-ratio-type statistics for testing efficiency is studied under both the null and the alternative hypotheses. The relevant nuisance parameter structure is established and finitesample bound procedures are proposed, which extend and improve available Gaussian-specific bounds. Finite-sample distributional invariance results are also demonstrated analytically for the HAC statistic proposed by MacKinlay and Richardson (1991). We study invariance to portfolio repacking for the tests and confidence sets proposed. The statistical properties of the proposed methods are analysed through a Monte Carlo study and compared with alternative available methods. Empirical results on NYSE returns show that exact confidence sets are very different from asymptotic ones, and allowing for non-Gaussian distributions affects inference results. Simulation and empirical evidence suggests that likelihood-ratio-type statistics with p-values corrected using the Maximized Monte Carlo test method—are generally preferable to their multivariate Fieller-Hotelling-HAC counterparts from the viewpoints of size control and power.

*Key words*: capital asset pricing model, CAPM, Black, mean-variance efficiency, non-normality, weak identification, Fieller, multivariate linear regression, uniform linear hypothesis, exact test, Monte Carlo test, bootstrap, nuisance parameters, GARCH, portfolio repacking

JEL Codes: C12, C33, C15, G12, G14

#### 1. INTRODUCTION

One of the most important extensions of the Capital Asset Pricing Model (CAPM) consists in allowing for the absence of a risk-free asset. From a theoretical viewpoint, this can be due to

restrictions on borrowing (Black, 1972) or an investor's "riskless" borrowing rate that exceeds the Treasury bill rate (Brennan, 1971). In this case, portfolio mean-variance efficiency is defined using the expected return in excess of the zero-beta portfolio. The latter is however unobservable which leads to considerable empirical difficulties.

Indeed, there are two basic approaches to estimating and assessing this version of the CAPM (denoted below as BCAPM). The first one uses a "two-pass" approach that may be traced back to Black *et al.* (1972) and Fama and MacBeth (1973): *betas* are first estimated from time series regressions for each security, and then the zero-beta rate is estimated by a cross-sectional regression on these *betas*. This raises errors-in-variables problems that affect statistical inference in both finite and large samples. The second approach—which originates in the work of Jensen (1968)—avoids this problem by using a multivariate linear regression as the basic statistical framework. In this article, we focus on the latter approach and consider two basic problems: (1) testing portfolio efficiency; (2) building a reliable confidence set for the zero-beta rate.

For clarity, let  $R_{it}$ ,  $i=1,\ldots,n$ , be the returns on n securities in period t, and  $\tilde{R}_{Mt}$  the return on a market benchmark for  $t=1,\ldots,T$ , and consider the n equations  $(i=1,\ldots,n)$  associated with the time series regressions of  $R_{it}$  on a constant and  $\tilde{R}_{Mt}$ , where the individual-equation disturbances are heteroskedastic and contemporaneously cross-correlated; let  $\Sigma = K'K$  refer to the error scale (or variance/covariance) matrix. If the intercepts from these n equations (the alphas) are denoted  $a_i$ , and the coefficients on the benchmark regressor (the betas) are denoted  $\beta_i$ ,  $i=1,\ldots,n$ , then the BCAPM equilibrium relations imply the following: there is a scalar  $\gamma$ , the return on the zero-beta portfolio, such that  $a_i - \gamma(1 - \beta_i) = 0$ ,  $i=1,\ldots,n$ . Our aim consists in assessing these constraints (denoted below as  $\mathcal{H}_B$ ) as well as estimating  $\gamma$ .

The above-cited literature provides analytical formulae for Gaussian likelihood-ratio statistics, the maximum likelihood estimator (MLE) of  $\gamma$  (denoted below as  $\hat{\gamma}$ ), and for a conformable asymptotic variance estimator [denoted Var( $\hat{\gamma}$ )]. It is however difficult to find reliable critical points in this context. While Gibbons (1982) used an asymptotic  $\chi^2$  critical value for the likelihood-ratio statistic, subsequent authors found this could lead to serious over-rejections, so various finite-sample corrections—such as bounds—have been suggested; see Shanken (1985, 1986, 1996), Stewart (1997), Zhou (1991, 1995), and Velu and Zhou (1999). These corrections depend crucially on normality, which may be inappropriate for financial data (see Fama, 1965; Richardson and Smith, 1993; Dufour *et al.*, 2003 and Beaulieu *et al.*, 2005, 2007, 2009, 2010). Furthermore, evidence on the properties of the confidence interval based on Var( $\hat{\gamma}$ ) is unavailable. Despite the simplicity of the above framework, discrepancies between asymptotic and finite-sample distributions are not surprising. Indeed, three difficulties deserve notice.

(1) Dimensionality: as n increases, the dimension of the scale matrix  $\Sigma$  grows rapidly and available degrees-of-freedom decrease conformably.<sup>3</sup> Even in linear or standard setups where the relevant asymptotic distributions may be free of  $\Sigma$ , this matrix can still affect the distributions in finite samples. Furthermore, positive definite estimates of  $\Sigma$  require a large T relative to n, so portfolios rather than securities are often used in practice.

<sup>1.</sup> See Litzenberger and Ramaswamy (1979), Banz (1981), Roll (1985), Chen et al. (1986), Shanken (1992), Kim (1995), Shanken and Zhou (2007), Lewellen et al. (2009), Kan et al. (2012), and Kleibergen (2009).

<sup>2.</sup> For other work based on the multivariate regression approach to CAPM analysis, see Gibbons (1982), Jobson and Korkie (1982), Kandel (1984, 1986), Amsler and Schmidt (1985), Shanken (1985, 1986, 1996), Kandel and Stambaugh (1989), Zhou (1991), Shanken (1992), Fama and French (1993), Chou (2000), Fama and French (2004), and Perold (2004).

<sup>3.</sup> See Shanken (1996), Campbell *et al.* (1997), Dufour and Khalaf (2002), Beaulieu *et al.* (2005, 2007, 2009, 2010), Sentana (2009), and the references therein.

- (2) Portfolio repacking (see Kandel and Stambaugh, 1989): to preserve meaningful pricing relations when portfolios are used, transformations of the return vector  $R_t = (R_{1t}, ..., R_{nt})'$  into  $R_t^* = AR_t$  where A is an  $n \times n$  invertible matrix such that  $A\iota_n = \iota_n$  and  $\iota_n$  is an n-dimensional vector of ones, should ideally not affect inference.
- (3) *Identification*: as  $\beta_i \to 1$ ,  $\gamma$  becomes weakly identified. Weak identification (WI) strongly affects the distributions of estimators and test statistics, leading to unreliable inference even asymptotically.<sup>4</sup> This should not be taken lightly: reported *betas* are often close to one (see *e.g.* Fama and MacBeth, 1973). Further, even if estimated *betas* are not close to one, irregularities associated with WI are not at all precluded [in view of (1) and (2) above]. Indeed, in the regression of  $R_t^*$  [from (2)] on a constant and  $\tilde{R}_{Mt}$ , with intercepts  $a_i^*$  and slopes  $\beta_i^*$ ,  $a_i \gamma(1 \beta_i) = 0$ ,  $i = 1, \ldots, n$ , for any  $\gamma$  and A. Portfolio repacking alters *betas* along with scale yet preserves the definition of  $\gamma$ , leading to identification problems as  $\beta_i^* \to 1$ . So the *betas* and scale parameters play a role in identifying  $\gamma$ .

Our aim in this article consists in providing inference methods that are robust to dimensionality and identification problems, whose outcomes are invariant to portfolio repacking. We first consider the problem of estimating  $\gamma$ . We show by simulation that available procedures provide poor coverage. So we propose exact confidence sets based on "inverting" exact tests for specific values of  $\gamma$ , *i.e.* the set of values not rejected by these tests. This method is a generalization of the classical procedure proposed by Fieller (1954) to estimate parameter ratios.<sup>5</sup>

To introduce the Fieller-type method in its simplest form with reference to the problem at hand, suppose (for illustrative purposes) that we aim at estimating  $\gamma$  from the univariate regression of the return of the i-th security ( $R_{it}$ ) on a constant and  $\tilde{R}_{Mt}$ , so that  $\gamma = -a_i/\delta_i$  where  $\delta_i = (\beta_i - 1)$ . Let  $\hat{a}_i$  and  $\hat{\delta}_i$  denote the OLS estimates from this regression, with estimated variances and covariance  $\text{Var}(\hat{a}_i)$ ,  $\text{Var}(\hat{\delta}_i)$  and  $\text{Cov}(\hat{a}_i,\hat{\delta}_i)$ . For each possible value  $\gamma_0$  of the ratio, consider the t-statistic  $t_i(\gamma_0) = (\hat{a}_i + \gamma_0 \hat{\delta}_i)/[\text{Var}(\hat{a}_i) + \delta_0^2 \text{Var}(\hat{\delta}_i) + 2\delta_0 \text{Cov}(\hat{a}_i,\hat{\delta}_i)]^{1/2}$  for testing  $\mathcal{H}_i(\gamma_0): a_i + \gamma_0 \delta_i = 0$ . Then, we obtain a confidence set with level  $1 - \alpha$  for  $\gamma$  by finding the set of  $\gamma_0$  values that are not rejected at level  $\alpha$  using  $t_i(\gamma_0)$  and a standard (normal or Student-t) two-tailed critical value  $z_{\alpha/2}$ . This means that we collect all  $\gamma_0$  values such that  $|t_i(\gamma_0)| \leq z_{\alpha/2}$  or alternatively such that  $(\hat{a}_i + \gamma_0 \hat{\delta}_i)^2 \leq z_{\alpha/2}^2(\text{Var}(\hat{a}_i) + \gamma_0^2 \text{Var}(\hat{\delta}_i) + 2\gamma_0 \text{Cov}(\hat{a}_i,\hat{\delta}_i))$ , leading to a second degree inequality in  $\gamma_0$ . The resulting confidence set has level  $1 - \alpha$  irrespective whether  $\delta_i$  is zero or not. In this article, we generalize this method to account for the multivariate setup where  $\gamma$  appears, allowing for Gaussian and non-Gaussian error distributions, as well as conditional heteroskedasticity. Empirically, we focus on multivariate Student-t and normal mixture distributions, as well as Gaussian GARCH.

To do so, we consider two statistics [denoted  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$ ] for testing  $\mathcal{H}(\gamma_0)$ :  $a_i + \gamma_0 \delta_i = 0$ ,  $i = 1, \ldots, n$ .  $LR(\gamma_0)$  is the likelihood-ratio statistic for the Gaussian error model, while  $\mathcal{J}(\gamma_0)$  is a multivariate Fieller-type statistic based on a generalized Hotelling statistic to which a correction for heteroskedasticity and autocorrelation (HAC) has been applied (as in MacKinlay and Richardson, 1991; Ravikumar *et al.*, 2000; and Ray and Savin, 2008). Using any one of these tests, we can build confidence sets by finding the values of  $\gamma_0$  that are not rejected at level  $\alpha$ . This requires a distributional theory for the test statistics. While an F-based cut-off point is

<sup>4.</sup> See, e.g. Dufour (1997, 2003), Staiger and Stock (1997), Wang and Zivot (1998), Zivot et al. (1998), Dufour and Jasiak (2001), Kleibergen (2002, 2005, 2009), Stock et al. (2002), Moreira (2003), Dufour and Taamouti (2005, 2007), and Andrews et al. (2006).

<sup>5.</sup> For the ratio of the means of two normal variables with equal variances, Fieller gave a solution that avoids non-regularities arising from a close-to-zero denominator. Extensions to univariate regressions or to several ratios with equal denominators can be found in Zerbe (1978), Dufour (1997), and Bolduc *et al.* (2010).

available for  $LR(\gamma_0)$  in the i.i.d. Gaussian case (see Beaulieu *et al.*, 2007 and Gibbons *et al.*, 1989), we show in a simulation study that usual asymptotic critical points perform poorly especially for  $\mathcal{J}(\gamma_0)$ .

To deal with such difficulties, we apply the maximized Monte Carlo (MMC) test procedure (Dufour, 2006) to obtain finite-sample p-values for  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$  in models with non-Gaussian and/or non-i.i.d. errors, as follows: a (simulated) p-value function conditional on relevant nuisance parameters is numerically maximized (with respect to these parameters), and the test is significant at level  $\alpha$  if the largest p-value is not larger than  $\alpha$ . The parametric bootstrap relates to the MMC method, in the sense that the maximization step is replaced by a unique p-value estimation, based on a consistent nuisance parameter estimate. For the GARCH case, such estimates may be unreliable in high-dimensional models; we show that the MMC method avoids this problem, with minimal power costs.

To implement this approach efficiently, it is important to characterize the nuisance parameters in the null distributions of the test statistics. We show that the null distributions of both  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$  do not depend on B and  $\Sigma$ , so the only nuisance parameters characterize the form of error distribution: the degree of freedom for Student-t distributions, the mixing probability and scaleratio parameters for normal mixtures, or GARCH parameters. While related invariance results are available for specific test problems using the multivariate regression likelihood ratio statistic [see Dufour and Khalaf (2002) and the references therein], our results on the HAC statistic  $\mathcal{J}(\gamma_0)$  are new to the literature and underscore the usefulness of MacKinlay and Richardson's (1991) statistic.

Because an F-based exact cut-off is available for the Gaussian case, we show that the confidence set which inverts  $LR(\gamma_0)$  can be obtained by solving a quadratic inequation. For non-i.i.d. or non-Gaussian distributions, we implement a numerical search running the MMC method for each choice for  $\gamma_0$ . Inverting  $\mathcal{J}(\gamma_0)$  requires numerical methods even for the Gaussian case. Furthermore, we show that all proposed confidence sets provide relevant information on whether efficiency is supported by the data, a property not shared by standard confidence intervals. Indeed, our confidence sets may turn out to be empty, which occurs when all possible values of  $\gamma$  are rejected.

We next consider testing efficiency in the BCAPM context. We study likelihood-ratio and multivariate Fieller-HAC criteria based on minimizing (over  $\gamma_0$ ) the above defined  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$  statistics. We show that the exact distribution of  $\min_{\gamma_0} \{LR(\gamma_0)\}$  depends on a reduced number of nuisance parameters that are functions of both B and  $\Sigma$ . We also generalize Shanken, 1986 exact bound test beyond the Gaussian model, and propose a tighter bound, which involves a numerical search for the tightest cut-off point, based on the MMC method. The MMC-based bound is also extended to the  $\min_{\gamma_0} \{ \mathcal{J}(\gamma_0) \}$  case. This approach, in conjunction with the above-defined confidence set based on  $\mathcal{J}(\gamma_0)$ , provides an interesting alternative to available GMM estimation methods [including the case recently analysed by Shanken and Zhou (2007)].

We present a simulation study to document the properties of the proposed procedures relative to available alternatives. In particular, we contrast problems arising from small samples with those caused by fundamentally flawed asymptotic approximations. We next examine efficiency of the market portfolio for monthly returns on New York Stock Exchange (NYSE) portfolios, built from the University of Chicago Center for Research in Security Prices (CRSP) 1926–1995 data base. We find more support for efficiency under the non-normal or non-i.i.d. hypothesis. Exact confidence sets for  $\gamma$  considerably differ from asymptotic ones, and the Fieller-HAC confidence sets are much wider than the GARCH corrected likelihood-ratio-based ones.

<sup>6.</sup> This procedure is based on the following fundamental property: when the distribution of a test statistic depends on nuisance parameters, the desired level  $\alpha$  is achieved by comparing the largest *p*-value (over all nuisance parameters consistent with the null hypothesis) with  $\alpha$ .

This article is organized as follows. Section 2 sets the framework and discusses identification of  $\gamma$ . In Section 3, we propose finite-sample tests for specific values of  $\gamma$ , and the corresponding exact confidence set are derived in Section 4. The exact distribution of the likelihood-ratio efficiency test statistic is established in Section 5, and bound procedures are proposed in Section 6. The simulation study is reported in Section 7. Our empirical analysis is presented in Section 8. We conclude in Section 9.

#### 2. MODEL AND ZERO-BETA IDENTIFICATION

Let  $R_{it}$ , i = 1, ..., n, be the returns on n securities in period t, and  $\tilde{R}_{Mt}$  the return on a market benchmark (t = 1, ..., T). Our analysis of the BCAPM model is based on the following standard multivariate regression setup (Gibbons, 1982; Shanken, 1986; MacKinlay, 1987):

$$R_{it} - \tilde{R}_{Mt} = a_i + (\beta_i - 1)\tilde{R}_{Mt} + u_{it}, \quad i = 1, ..., n, t = 1, ..., T,$$
 (2.1)

where  $u_{it}$  is a random disturbance. The testable implication of the BCAPM on (2.1) is the following one: there is a scalar  $\gamma$ , the return on the zero-beta portfolio, such that

$$\mathcal{H}_{\mathbf{R}}: a_i + \gamma \delta_i = 0, \quad \delta_i = \beta_i - 1, \quad i = 1, \dots, n, \quad \text{for some } \gamma \in \Gamma,$$
 (2.2)

where  $\Gamma$  is the set of "admissible" values for  $\gamma$ . Since  $\gamma$  is unknown,  $\mathcal{H}_B$  is non-linear. The latter can be viewed as the union of more restrictive linear hypotheses of the form

$$\mathcal{H}(\gamma_0): a_i + \gamma_0 \delta_i = 0, i = 1, \dots, n, \tag{2.3}$$

where  $y_0$  is specified. This observation underlies the exact inference approach proposed here.

## 2.1. Notation

The above model is a special case of the following multivariate regression:

$$Y = XB + U \tag{2.4}$$

where  $Y = [Y_1, ..., Y_n]$  is  $T \times n$ , X is  $T \times k$  of rank k,  $U = [U_1, ..., U_n] = [V_1, ..., V_T]'$ . For (2.1),  $Y = [R_1, ..., R_n]$ ,  $X = [\iota_T, \tilde{R}_M]$ ,  $R_i = (R_{i1}, ..., R_{iT})'$ ,  $\tilde{R}_M = (\tilde{R}_{M1}, ..., \tilde{R}_{MT})'$ ,  $B = [a, \beta]'$ ,  $a = (a_1, ..., a_n)'$ ,  $\beta = (\beta_1, ..., \beta_n)'$ , and  $\iota_T$  refers to a T-dimensional vector of ones.

Throughout the article, we use the following notation.  $P_{(B,K)}$  represents the distribution of Y when the parameters are (B,K), where K represents parameters not included in B (such as parameters of the error distribution). For any matrix  $n \times k$  matrix A,  $M(A) = I - A(A'A)^{-}A'$  and vec(A) is the  $(nk) \times 1$  vector obtained by stacking the columns of A on top of each other. We also use the following equivalent forms for the model, parameters, and hypotheses considered:

$$\tilde{Y} = Y - \tilde{R}_{\mathbf{M}} \iota_n' = XC + U, \tag{2.5}$$

where 
$$C = B - \Delta = [a, \beta - \iota_n]', \quad \Delta = [0, \iota_n]';$$
 (2.6)

$$\tilde{\mathcal{H}}(\gamma_0): H(\gamma_0)C = 0$$
, for  $\gamma_0$  specified, (2.7)

$$\widetilde{\mathcal{H}}(\gamma_0)$$
:  $R(\gamma_0)\vartheta = 0$ , for  $\gamma_0$  specified, (2.8)

where 
$$H(\gamma_0) = (1, \gamma_0)$$
,  $R(\gamma_0) = H(\gamma_0) \otimes I_n$ ,  $\vartheta = \text{vec}(C')$ ,  $R(\gamma)\vartheta = [H(\gamma)C]'$ , (2.9)

$$H(\gamma)C = 0 \Leftrightarrow R(\gamma)\vartheta = 0$$
, for all  $\gamma$ ; (2.10)

$$\widetilde{\mathcal{H}}_{B}: H(\gamma)C = 0$$
, for some  $\gamma \in \Gamma$ , (2.11)

$$\widetilde{\mathcal{H}}_{B}: R(\gamma)\vartheta = 0$$
, for some  $\gamma \in \Gamma$ . (2.12)

Note that (2.7) and (2.8) represent equivalent representations of  $\tilde{\mathcal{H}}(\gamma_0)$ , and similarly (2.11) and (2.12) for  $\tilde{\mathcal{H}}_B$ .

## 2.2. Distributional assumptions

We further assume that we can condition on  $\tilde{R}_{\rm M}$  and

$$V_t = (u_{1t}, \dots, u_{nt})' = K'W_t, \quad t = 1, \dots, T, \quad W_t = (W_{1t}, \dots, W_{nt})',$$
 (2.13)

where K is unknown and non-singular,  $W = [W_1, ..., W_T]'$  is independent of X, and the distribution of W is either fully specified or determined up to an unknown distributional shape parameter  $\nu$ . (2.13) can also be rewritten in matrix form as U = WK.

We first present results that require no further regularity assumptions. We use additional restrictions, which entail that the distribution of W belongs to a specific family  $\mathcal{H}_W(\mathcal{D}, \nu)$ , where  $\mathcal{D}$  represents a distribution type and  $\nu \in \Omega_{\mathcal{D}}$  a nuisance parameter characterizing the distribution. In particular, we consider the multivariate normal  $(\mathcal{D}_N)$ , Student- $t(\mathcal{D}_t)$ , and normal mixture  $(\mathcal{D}_m)$  distributions:

$$\mathscr{H}_{W}(\mathscr{D}_{N}): W_{t} \overset{i.i.d}{\sim} N[0, I_{n}],$$
 (2.14)

$$\mathscr{H}_{W}(\mathscr{D}_{t},\kappa):W_{t}=Z_{1t}/(Z_{2t}/\kappa)^{1/2}, Z_{1t}\overset{i.i.d}{\sim} N[0,I_{n}], Z_{2t}\overset{i.i.d}{\sim} \chi^{2}(\kappa),$$
 (2.15)

$$\mathcal{H}_{W}(\mathcal{D}_{m}, \pi, \omega): W_{t} = I_{t}(\pi) Z_{1t} + [1 - I_{t}(\pi)] Z_{3t}, Z_{3t} \stackrel{i.i.d}{\sim} N[0, \omega I_{n}], 0 < \pi < 1,$$
(2.16)

where  $Z_{2t}$  and  $Z_{3t}$  are independent of  $Z_{1t}$ , and  $I_t(\pi)$  is an indicator random variable independent of  $(Z_{1t}, Z_{3t})$  such that  $P[I_t(\pi) = 0] = 1 - P[I_t(\pi) = 1] = \pi$ . So, in (2.13),  $v = \kappa$  under (2.15), and  $v = (\pi, \omega)$  under (2.16). If  $E(W_t W_t') = I_n$ , the covariance matrix of  $V_t$  is  $\Sigma = K'K$ , so  $\Sigma$  is positive definite without further restrictions. If K' is lower triangular, K' and K correspond to the Cholesky factors of  $\Sigma$ .

Time-dependence may be taken into account by an appropriate specification of the distribution of  $W_t$ ,  $t=1, \ldots, T$ . Since time-varying volatility is prevalent in financial data, we consider the parametric GARCH structure:

$$u_{it} = w_{it} h_{it}^{\frac{1}{2}}, \qquad h_{it} = (1 - \phi_{1i} - \phi_{2i}) \sigma_i^2 + \phi_{1i} w_{i,t-1}^2 + \phi_{2i} h_{i,t-1},$$
 (2.17)

where  $w_{it}$  are uncorrelated standard normal variables. This process may easily be reparameterized as in (2.13), where K is a diagonal matrix with diagonal terms  $(1-\phi_{1i}-\phi_{2i})^{1/2}\sigma_i$ ,  $i=1,\ldots,n$ , and each  $W_{it}$  follows a univariate stationary GARCH process with unit intercept. Conforming with the above notation, we refer to this distributional hypothesis as  $\mathscr{H}_W(\mathscr{D}_G,\phi)$ , where  $\phi$  is the  $2n \times 1$  vector  $(\phi_{11},\ldots,\phi_{1n},\phi_{21},\ldots,\phi_{2n})^7$ 

<sup>7.</sup> Ideally, a multivariate GARCH structure may be considered if T is sufficiently large relative to n; see Bauwens *et al.* (2006) for a recent survey. We adopt (2.17) since our empirical analysis relies on monthly data with 12 portfolios over 5 year subperiods (*i.e.* T = 60 and n = 12).

## 2.3. Weak identification

Even though  $a_i$  and  $\beta_i$  are identifiable,  $\gamma$  is defined through a non-linear transformation that may fail to be well defined: the ratio  $\gamma = a_i/(1-\beta_i)$  is not defined or, equivalently, the equation  $a_i = \gamma(1-\beta_i)$  does not have a unique solution when  $\beta_i = 1$ .

In such situations, the distributions of many usual test statistics become non-standard, so the corresponding tests are unreliable and the associated confidence sets invalid. In particular, asymptotic standard errors are unreliable measures of uncertainty, and standard asymptotically justified *t*-type tests and confidence intervals have sizes that may deviate arbitrarily from their nominal levels; see the literature on weak identification [as reviewed, for example, in Dufour (2003) and Stock *et al.* (2002)]. Both the finite and large-sample distribution theory of most test statistics can be affected.

While it is straightforward to see that  $\beta_i = 1$  corresponds to a discontinuity, the analysis below reveals this is not the whole story. In particular, we study the properties of estimators and test statistics following data transformations of the form  $\tilde{Y}_* = \tilde{Y}A$ , where A is any non-singular fixed matrix of order n. On comparing (2.1) to its transformed counterpart, we see that irregularities cannot be safely assumed away, even when observed *betas* are not close to one.

#### 2.4. Standard estimators and test statistics

One of the most common inference methods in this context relies on the log-likelihood

$$\ln[L(Y,B,\Sigma)] = -\frac{nT}{2}(2\pi) - \frac{T}{2}\ln(|\Sigma|) - \frac{1}{2}\text{tr}[\Sigma^{-1}(Y - XB)'(Y - XB)].$$
(2.18)

The unrestricted MLE of B and  $\Sigma$  are:

$$\hat{B} = (X'X)^{-1}X'Y = [\hat{a}, \hat{\beta}]', \quad \hat{\Sigma} = \hat{U}'\hat{U}/T,$$

where  $\hat{U} = Y - X\hat{B}$ ,  $\hat{a} = (\hat{a}_1, \dots, \hat{a}_n)'$  and  $\hat{\beta} = (\hat{\beta}_1, \dots, \hat{\beta}_n)'$ . If  $\hat{C}$  is the MLE of C in (2.5), the corresponding estimate of  $\theta = \text{vec}(C')$  is

$$\hat{\vartheta} = \text{vec}(\hat{C}') \tag{2.19}$$

where  $\vartheta$  is defined in (2.6).

**2.4.1. Gaussian-based statistics.** The likelihood-ratio statistic to test  $\mathcal{H}(\gamma_0)$  where  $\hat{\Sigma}(\gamma_0)$  is the MLE of  $\Sigma$  under  $\mathcal{H}(\gamma_0)$  is:

$$LR(\gamma_0) = T \ln[\Lambda(\gamma_0)], \quad \Lambda(\gamma_0) = |\hat{\Sigma}(\gamma_0)|/|\hat{\Sigma}| = \frac{n}{T - n - 1} \mathscr{F}(\gamma_0) + 1,$$
 (2.20)

$$\hat{\Sigma}(\gamma_0) = \hat{\Sigma} + \frac{1}{T} \hat{B}' H(\gamma_0)' [H(\gamma_0)(X'X)^{-1} H(\gamma_0)']^{-1} H(\gamma_0) \hat{B}, \qquad (2.21)$$

$$\mathscr{F}(\gamma_0) = \frac{T - n - 1}{n} \frac{\left(\hat{a} + \hat{\delta}\gamma_0\right)' \hat{\Sigma}^{-1} \left(\hat{a} + \hat{\delta}\gamma_0\right)}{1 + \left[\left(\hat{\mu}_M - \gamma_0\right)^2 / \hat{\sigma}_M^2\right]},\tag{2.22}$$

$$\hat{\mu}_{M} = \frac{1}{T} \sum_{t=1}^{T} \tilde{R}_{Mt}, \quad \hat{\sigma}_{M}^{2} = \frac{1}{T} \sum_{t=1}^{T} (\tilde{R}_{Mt} - \hat{\mu}_{M})^{2}, \quad \hat{\delta} = \hat{\beta} - \iota_{n}.$$
 (2.23)

 $\mathscr{F}(\gamma_0)$  is the Hotelling statistic for testing  $\mathscr{H}(\gamma_0)$ . Even though this statistic may look like a "Wald-type" statistic, the "covariance matrix" used for  $\hat{a} + \hat{\delta}\gamma_0$ —namely,

 $(1+[(\hat{\mu}_M-\gamma_0)^2/\hat{\sigma}_M^2])\hat{\Sigma}$  – depends on  $\gamma_0$  (the tested value of  $\gamma$ ). This characteristic makes it similar to Fieller-type statistics in univariate linear regressions (see Dufour, 1997), so  $\mathcal{F}(\gamma_0)$  can be described as a "Fieller-Hotelling-type statistic". We will see below that this feature plays a crucial role in avoiding distributional problems associated with weak identification.

The likelihood-ratio criterion to test  $\mathcal{H}_{B}$  is

$$LR_{\rm B} = T\ln(\Lambda_{\rm B}) = \inf\{LR(\gamma_0): \gamma_0 \in \Gamma\} = LR(\hat{\gamma}), \tag{2.24}$$

$$\Lambda_{\mathbf{B}} = |\hat{\Sigma}_{\mathbf{B}}|/|\hat{\Sigma}|, \quad |\hat{\Sigma}_{\mathbf{B}}| = \inf\{|\hat{\Sigma}(\gamma_0)| : \gamma_0 \in \Gamma\},\tag{2.25}$$

where  $\hat{\Sigma}_{B}$  is the MLE of  $\Sigma$  under  $\mathcal{H}_{B}$  and  $\hat{\gamma}$  is the unrestricted MLE of  $\gamma$ ; see Shanken (1986). The log-likelihood for (2.5) is

$$\ln\left[\tilde{L}(\tilde{Y},C,\Sigma)\right] = \ln\left[L(Y-\tilde{R}_{\mathrm{M}}\iota_{n}',B-\Delta,\Sigma)\right] = \ln\left[L(Y,B,\Sigma)\right] \tag{2.26}$$

and the likelihood-ratio statistics for testing  $\tilde{\mathcal{H}}(\gamma_0)$  and  $\tilde{\mathcal{H}}_B$  coincide with  $LR(\gamma_0)$  and  $LR_B$ . Throughout the article, we treat  $LR_B$  and  $LR(\gamma_0)$  as quasi likelihood-ratio (QLR) criteria and the associated MLEs as quasi maximum likelihood (QML) estimators. We denote the observed value of these statistics as  $LR_{\rm B}^{(0)}$  and  $LR^{(0)}(\gamma_0)$ , respectively. A Wald-type formula for an asymptotic information-matrix-based standard error associated

with  $\hat{\gamma}$  is provided by (Campbell *et al.*, 1997, Chapter 5, equation 5.3.81):

$$\operatorname{Var}(\hat{\gamma}) = \frac{1}{T} \left[ 1 + \frac{(\hat{\mu}_M - \gamma)^2}{\hat{\sigma}_M^2} \right] [(\iota_n - \beta)' \Sigma^{-1} (\iota_n - \beta)]^{-1}.$$
 (2.27)

Whereas corrections may be derived for the non-Gaussian case [as in Barone-Adesi et al. (2004) who study a related asset pricing problem], general results on inference in the presence of identification failure (or weak identification) indicates that a "variance" estimator that does not depend on the tested value  $\gamma_0$ —which is the case for  $Var(\hat{\gamma})$ —cannot lead to valid pivotal functions, so the associated "asymptotic" confidence sets and tests are fundamentally invalid; see Dufour (1997). Here, for example, it is easy to see that the above formula cannot be valid when  $\beta = \iota_n$ , and problematic when  $\beta \subseteq \iota_n$ .

**2.4.2.** Fieller-Hotelling-HAC statistics.  $\mathscr{F}(\gamma_0)$  may be viewed as a Hotelling-type statistic based on the standardized distance between  $\hat{a}+\hat{\delta}\gamma_0$  and zero, which conveys an asymptotic least-squares (Gouriéroux et al., 1985; Gouriéroux and Monfort, 1995, Ch. 9) and a GMM interpretation of  $\hat{\gamma}$ . This may be exploited to allow for serial dependence, for example by using a properly corrected weighting matrix, as done for example by MacKinlay and Richardson (1991), Ravikumar et al. (2000), and Ray and Savin (2008). This suggests the following statistic:

$$\mathscr{J}(\gamma_0) = T \hat{\vartheta}' \mathsf{R}(\gamma_0)' \left\{ \mathsf{R}(\gamma_0) \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] \mathsf{S}_T \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] \mathsf{R}(\gamma_0)' \right\}^{-1} \mathsf{R}(\gamma_0) \hat{\vartheta}, \quad (2.28)$$

where  $R(\gamma_0)$  is defined in (2.9),

$$S_T = \Psi_{0,T} + \sum_{j=1}^{q} \left( \frac{q-j}{q} \right) \left[ \Psi_{j,T} + \Psi'_{j,T} \right], \quad \Psi_{j,T} = \frac{1}{T} \sum_{t=j+1}^{T} \left( X_t \otimes \hat{U}_t \right) \left( X_{t-j} \otimes \hat{U}_{t-j} \right)',$$

and  $\hat{U}_t'$  is the *t*-th row of  $\hat{U}$ . A crucial feature of  $\mathcal{J}(\gamma_0)$  comes from the fact that the covariance matrix estimator used for  $R(\gamma_0)\hat{\vartheta}$  depends on  $\gamma_0$  (the tested value of  $\gamma$ ). Since it is a HAC-modified version of  $\mathcal{F}(\gamma_0)$ ,  $\mathcal{J}(\gamma_0)$  can be described as a "Fieller-Hotelling-HAC" statistic or " $\mathcal{F}(\gamma_0)$ -HAC" statistic.

Under  $\mathcal{H}(\gamma_0)$ ,  $\mathcal{J}(\gamma_0)$  follows a  $\chi^2(n)$  distribution asymptotically. This result does not require any assumption on the identification of  $\gamma$ . A GMM estimator  $\tilde{\gamma}$  of  $\gamma$  can be obtained by solving the problem

$$\mathcal{J}_{B} = \inf \left\{ \mathcal{J}(\gamma_{0}) : \gamma_{0} \in \Gamma \right\} = \mathcal{J}(\tilde{\gamma}). \tag{2.29}$$

We denote the observed value of these statistics as  $\mathcal{J}_{B}^{(0)}$  and  $\mathcal{J}^{(0)}(\gamma_{0})$ , respectively.

# 3. IDENTIFICATION-ROBUST MONTE CARLO TESTS FOR $\gamma$

We will now derive the exact null distribution of  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$  under  $\mathcal{H}(\gamma_0)$ , where  $\gamma_0$  is known. We then show how this result can be used to obtain exact Monte Carlo p-values. This will allow us to build confidence sets for  $\gamma$  and yield a way of testing efficiency.

# 3.1. Distribution of $LR(\gamma_0)$ and $\mathcal{J}(\gamma_0)$ under $\mathcal{H}(\gamma_0)$

Theorems 3.1 and 3.2 show that the null distribution of both  $LR(\gamma_0)$  and  $\mathcal{J}(\gamma_0)$  given X, is completely determined by X and the distribution of W given X. Proofs are given in the Appendix.

**Theorem 3.1.** DISTRIBUTION OF THE MEAN-VARIANCE CAPM TEST FOR A KNOWN ZERO-BETA RATE. Under (2.1), (2.13), and  $\mathcal{H}(\gamma_0)$ ,  $LR(\gamma_0)$  is distributed like

$$\overline{LR}(\gamma_0, W) = T \ln(\left| W' \overline{M}(\gamma_0) W \right| / \left| W' M W \right|) \tag{3.1}$$

where  $\bar{M}(\gamma_0) = M(X) + X(X'X)^{-1}H(\gamma_0)'[H(\gamma_0)(X'X)^{-1}H(\gamma_0)']^{-1}H(\gamma_0)(X'X)^{-1}X'$ .

In the i.i.d. Gaussian case (2.14), we have:

$$[(T-1-n)/n][\Lambda(\gamma_0)-1] \sim F(n, T-1-n); \tag{3.2}$$

see Dufour and Khalaf (2002). This result was used by Gibbons *et al.* (1989) in studying efficiency with an observable risk-free rate. Indeed, testing  $\mathcal{H}(\gamma_0)$  is equivalent to testing whether the intercepts are jointly zero in a market model with returns in excess of  $\gamma_0$ .

**Theorem 3.2.** DISTRIBUTION OF THE FIELLER-HOTELLING-HAC STATISTIC FOR  $\gamma$ . *Under* (2.1), (2.13), and  $\mathcal{H}(\gamma_0)$ , the statistic  $\mathcal{J}(\gamma_0)$  defined in (2.28) is distributed like

$$\mathcal{J}(\gamma_0, W) = TH(\gamma_0)(X'X)^{-1}X'W\bar{\mathcal{Q}}(\gamma_0, W)^{-1}W'X(X'X)^{-1}H(\gamma_0)', \tag{3.3}$$

where

$$\begin{split} \bar{\mathcal{Q}}(\gamma_0, W) = \left[ H(\gamma_0) \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] \hat{S}_T \left[ \left( \frac{X'X}{T} \right)^{-1} H(\gamma_0)' \otimes I_n \right], \\ \hat{S}_T = \hat{\Psi}_{0,T} + \sum_{j=1}^q \left( \frac{q-j}{q} \right) \left[ \hat{\Psi}_{j,T} + \hat{\Psi}_{j,T}' \right], \quad \hat{\Psi}_{j,T} = \frac{1}{T} \sum_{t=j+1}^T \left( X_t \otimes \hat{W}_t \right) \left( X_{t-j} \otimes \hat{W}_{t-j} \right)', \end{split}$$

and  $\hat{W}_{t}'$  is the t-th row of  $\hat{W} = M(X)W$ .

For non-Gaussian distributions compatible with (2.13) [including the GARCH case (2.17)], Theorem 3.1 shows that the exact distribution of both statistics, although non-standard, may easily be simulated once X, the distribution of W and  $\gamma_0$  [given by  $\mathcal{H}(\gamma_0)$ ] are set. So the Monte Carlo (MC) test method can be easily applied; see Dufour (2006). Of course, the exactness feature does not hold when the assumptions used do not hold. In particular, this is the case if serial dependence is present, although the test based on  $\mathcal{J}(\gamma_0)$  remains asymptotically valid in the usual way.

## 3.2. Monte Carlo test method

This section explains the MC and MMC methods for implementing hypothesis tests based on a test *S* statistic. We also set associated notation, given the following assumptions on *S*.

In the context of model (2.4), consider a null hypothesis  $H_*$  which consists of (possibly nonlinear) restrictions on B, and an associated test statistic S. We reject  $H_*$  when  $S \ge c$ . Under  $H_*$ , S is distributed like a function  $\bar{S}(\eta, W)$  of W and a vector of parameters  $\eta \in \Xi$  (as well as the known matrix X), where  $\Xi$  describes the restricted parameter space (typically of dimension lower than the number of parameters in B and K). Finally, given (2.13), we can simulate W, once v has been specified [e,g, as suggested by one of the assumptions (2.14)–(2.16)].

For notation simplicity, the dependence upon X is implicit through the definition of  $\bar{S}$ . Examples of  $\bar{S}(.)$  include (3.1) and (3.3) in which case  $\eta$  corresponds to  $\gamma_0$  which is set by the null hypothesis. The MC method assesses the rank of the observed value of S denoted [denoted  $S^{(0)}$ ], relative to a finite number N of simulated statistics [denoted  $S^{(1)}, ..., S^{(N)}$ ] drawn under the null hypothesis. Given the above general assumptions, we can apply the following algorithm.

A1 Given: (i) a value of  $\nu$ , (ii) N draws  $W^{(1)}, \ldots, W^{(N)}$  from the distribution of W, (iii) a value of  $\eta$ , and (iv) the test function  $\bar{S}(\eta, W)$ , compute  $\bar{S}(\eta, W)$  leading to the vector

$$\bar{S}_N(\eta, \nu) = \left[\bar{S}\left(\eta, W^{(1)}\right), \dots, \bar{S}\left(\eta, W^{(N)}\right)\right]'. \tag{3.4}$$

A2 Build the MC p-value function

$$p_N[S^{(0)}|\bar{S}_N(\eta,\nu)] = \frac{NG_N[S^{(0)};\bar{S}_N(\eta,\nu)] + 1}{N+1},$$
(3.5)

$$G_N[S^{(0)}; \bar{S}_N(\eta, \nu)] = \frac{1}{N} \sum_{j=1}^N I_{[0,\infty)}[\bar{S}(W^{(j)}, \eta) - S^{(0)}],$$
 (3.6)

where  $I_A[x] = 1$ , if  $x \in A$ , and  $I_A[x] = 0$ , if  $x \notin A$ . Set  $\hat{p}_N(S; \eta, \nu) = p_N[S^{(0)} | \bar{S}_N(\eta, \nu)]$ . A3 If  $\nu$  and  $\eta$  are set by  $H_*$ , the test which rejects  $H_*$  when

$$\hat{p}_N(S;\eta,\nu) \le \alpha \tag{3.7}$$

is a test with level  $\alpha$  for  $H_*$ ; if furthermore the distribution of S is continuous under  $H_*$  and  $\alpha(N+1)$  is an integer [as assumed in the cases considered in this article], this test has size  $\alpha$ : the probability of rejection under the null hypothesis is equal to  $\alpha$ , for finite T and N. If  $\nu$  or  $\eta$  is not set by the null hypothesis, then maximize  $p_N[S^{(0)}|\bar{S}_N(\eta,\nu)]$  over

all the  $(\nu, \eta)$  values compatible with  $H_*$ , and reject the latter if the maximal p-value is less than or equal to  $\alpha$ . For any  $A \subseteq \Xi$  and  $E \subseteq \Omega_{\mathscr{D}}$ , let

$$\hat{p}_N(S; \eta, E) = \sup \{ \hat{p}_N(S; \eta, \nu) : \nu \in E \}, \tag{3.8}$$

$$\hat{p}_N(S; A, \nu) = \sup\{\hat{p}_N(S; \eta, \nu) : \eta \in A\},\tag{3.9}$$

$$\hat{p}_N(S; A, E) = \sup\{\hat{p}_N(S; \eta, \nu) : \eta \in A, \nu \in E\}, \tag{3.10}$$

where, by convention,  $\hat{p}_N(S; A, \cdot) = 0$  if A is empty, and  $\hat{p}_N(S; \cdot, E) = 0$  if E is empty. Then the probability of rejection under  $H_*$  based on either: (i)  $\hat{p}_N(S; \eta, \Omega_{\mathscr{D}})$  when  $\eta$  is set by  $H_*$ , (ii)  $\hat{p}_N(S; \Xi, \nu)$  when  $\nu$  is set by  $H_*$ , or (iii)  $\hat{p}_N(S; \Xi, \Omega_{\mathscr{D}})$  when both  $\nu$  and  $\eta$  are not set by  $H_*$ , is itself not larger than  $\alpha$  for finite T and N; see Dufour (2006).

A test for which the probability of rejection under the null hypothesis is not larger than  $\alpha$  is sometimes described as a *conservative* test. We prefer to rely on the more precise and traditional distinction between a level-correct test and a size-correct test. A test is level-correct for  $\alpha$  (or exact at level  $\alpha$ ) if the probability of rejection under all distributions compatible with the null hypothesis is *less than or equal* to  $\alpha$  (for finite T). In contrast, a test is *size*-correct for  $\alpha$  (or has exact size  $\alpha$ ) if the supremum of the probability of rejection under all distributions compatible with the null hypothesis is equal to  $\alpha$  (for finite T). It is clear that a size-correct test is level-correct, but the converse may not hold. Both these definitions account for the common situation where the rejection probability under the null hypothesis may depend on nuisance parameters, and thus vary over the null hypothesis. We use the term exact when underlying probability statements hold for finite T and N, whether null rejection probabilities are equal to or not larger than the hypothesized cut-off. We follow the same notation in the case of a confidence set. A confidence set with exact size  $1-\alpha$  implies that the infimum of the set's coverage probabilities compatible with the considered model is equal to  $1-\alpha$ , for finite T and N. In contrast, a confidence set with exact level  $\alpha$  implies that the set's coverage probability is greater than or equal to  $1-\alpha$ , for finite T and N. This definition implies that inverting a size-correct test leads to a size-correct confidence set, while inverting a level-correct test leads to a level-correct confidence set.

When a Monte Carlo p-value is computed given a hypothesized distribution, and where in this case, the underlying test statistic converges to any distribution (say  $\bar{F}$ ) the associated Monte Carlo test will remain asymptotically valid under any set of weaker assumptions for which the statistic still converges to the same limiting distribution  $\bar{F}$ . The implications for the present article are as follows: our tests though exact with the considered distributions on W remain asymptotically valid in a semi-parametric context under the usual set of assumptions typically required for the CAPM to hold.

# 3.3. Monte Carlo tests based on $LR(\gamma_0)$ and $\mathcal{J}(\gamma_0)$

With respect to the algorithm in the previous section, in the case of  $LR(\gamma_0)$ ,  $S^{(0)} \equiv LR^{(0)}(\gamma_0)$ ,  $\eta \equiv \gamma_0$ , and using (3.1),  $\bar{S}(\eta, W^{(i)}) = \overline{LR}(\gamma_0, W^{(i)})$ . With  $\bar{S}_N(\eta, \nu) = \overline{LR}_N(\gamma_0, \nu) = [\overline{LR}(\gamma_0, W^{(1)}), ..., \overline{LR}(\gamma_0, W^{(N)})]'$ , and using the *p*-value function  $p_N[.]$  from (3.5), let

$$\hat{p}_N(LR;\gamma_0,\nu) \equiv p_N \left[ LR^{(0)}(\gamma_0) \middle| \overline{LR}_N(\gamma_0,\nu) \right]. \tag{3.11}$$

As a result of Theorem 3.1, we have, under  $\mathcal{H}(\gamma_0)$  in conjunction with  $\mathcal{H}_W(\mathcal{D}, \nu)$  and given the notation from (3.8) to (3.10):

$$P[\hat{p}_N(LR; \gamma_0, \nu_0) \le \alpha] = \alpha, \text{ when } \nu = \nu_0, \tag{3.12}$$

$$P[\hat{p}_N(LR; \gamma_0, \Omega_{\mathscr{D}}) \le \alpha] \le \alpha, \text{ when } \nu \text{ may be unknown.}$$
 (3.13)

Similarly, in the case of  $\mathcal{J}(\gamma_0)$ ,  $S^{(0)} \equiv \mathcal{J}^{(0)}(\gamma_0)$ ,  $\eta \equiv \gamma_0$ , and using (3.3),  $\bar{S}(\eta, W^{(i)}) = \bar{\mathcal{J}}(\gamma_0, W^{(i)})$ . With  $\bar{S}_N(\eta, \nu) = (\gamma_0, \nu) = [\bar{\mathcal{J}}(\gamma_0, W^{(1)}), \dots, \bar{\mathcal{J}}(\gamma_0, W^{(N)})]'$ , let

$$\hat{p}_N(\mathcal{J}; \gamma_0, \nu) \equiv p_N \left[ \mathcal{J}^{(0)}(\gamma_0) \middle| \bar{\mathcal{J}}_N(\gamma_0, \nu) \right]. \tag{3.14}$$

As a result of Theorem 3.2, we have, under  $\mathcal{H}(\gamma_0)$  in conjunction with  $\mathcal{H}_W(\mathcal{D}, \nu)$ :

$$P[\hat{p}_N(\mathcal{J}; \gamma_0, \nu_0) \le \alpha] = \alpha, \text{ when } \nu = \nu_0, \tag{3.15}$$

$$P[\hat{p}_N(\mathcal{J}; \gamma_0, \Omega_{\mathcal{D}}) \le \alpha] \le \alpha, \text{ when } \nu \text{ may be unknown.}$$
 (3.16)

#### 4. IDENTIFICATION-ROBUST CONFIDENCE SETS FOR $\gamma$

Under  $\mathcal{H}_B$ , the ratios  $a_i/(1-\beta_i)$ , 1, ..., n, are equal. This definition of  $\gamma$  leads to the classical problem of inference on ratios from Fieller (1954). The problem here is clearly more complex, so to extend Fieller's arguments, we use the above-defined tests of  $\mathcal{H}(\gamma_0)$ .

## 4.1. Hotelling-Fieller confidence sets: the i.i.d. Gaussian case

Consider the Gaussian model given by (2.1), (2.13), and (2.14). In this case, under  $\mathcal{H}_0(\gamma_0)$ ,  $\mathcal{F}(\gamma_0)$  follows a Fisher distribution F(n, T-n-1); see (3.2). Let  $F_{\alpha}$  denote the cut-off point for a test with level  $\alpha$  based on the F(n, T-n-1) distribution. Then

$$CF_{\gamma}(\alpha) = \{ \gamma_0 \in \Gamma : \mathscr{F}(\gamma_0) \le F_{\alpha} \}$$
 (4.1)

has level  $1-\alpha$  for  $\gamma$ , *i.e.* the probability that  $\gamma$  be covered by  $CF_{\gamma}(\alpha)$  is not smaller than  $1-\alpha$ . Indeed,  $P[\gamma \in CF_{\gamma}(\alpha)] = 1-\alpha$ , that is  $CF_{\gamma}(\alpha)$  is size-correct. On noting that  $\mathscr{F}(\gamma_0) \leq F_{\alpha}$  can be rewritten as

$$M_F(\gamma_0) - \frac{nF_\alpha}{T - n - 1} N_F(\gamma_0) \le 0,$$
 (4.2)

$$M_F(\gamma_0) = (\hat{a} + \hat{\delta}\gamma_0)'\hat{\Sigma}^{-1}(\hat{a} + \hat{\delta}\gamma_0) = (\hat{\delta}'\hat{\Sigma}^{-1}\hat{\delta})\gamma_0^2 + (2\hat{\delta}'\hat{\Sigma}^{-1}\hat{a})\gamma_0 + \hat{a}'\hat{\Sigma}^{-1}\hat{a}, \tag{4.3}$$

$$N_F(\gamma_0) = 1 + \frac{(\hat{\mu}_M - \gamma_0)^2}{\hat{\sigma}_M^2} = \frac{1}{\hat{\sigma}_M^2} \gamma_0^2 - \frac{2\hat{\mu}_m}{\hat{\sigma}_M^2} \gamma_0 + 1 + \frac{\hat{\mu}_M^2}{\hat{\sigma}_M^2}, \tag{4.4}$$

we see, after a few manipulations, that  $CF_{\nu}(\alpha)$  reduces to a simple quadratic inequation:

$$CF_{\gamma}(\alpha) = \{ \gamma_0 \in \Gamma : A\gamma_0^2 + B\gamma_0 + C \le 0 \}, \tag{4.5}$$

$$A = \hat{\delta}' \hat{\Sigma}^{-1} \hat{\delta} - \left(\frac{nF_{\alpha}}{T - n - 1}\right) \frac{1}{\hat{\sigma}_{M}^{2}}, \quad B = 2\left[\hat{\delta}' \hat{\Sigma}^{-1} \hat{a} + \left(\frac{nF_{\alpha}}{T - n - 1}\right) \frac{\hat{\mu}_{M}}{\hat{\sigma}_{M}^{2}}\right], \tag{4.6}$$

$$C = \hat{a}' \hat{\Sigma}^{-1} \hat{a} - \left(\frac{nF_{\alpha}}{T - n - 1}\right) \left[1 + \frac{\hat{\mu}_{M}^{2}}{\hat{\sigma}_{M}^{2}}\right]. \tag{4.7}$$

For  $\Gamma = \mathbb{R}$ , the resulting confidence set can take several forms depending on the roots of the polynomial  $A\gamma_0^2 + B\gamma_0 + C$ : (a) a closed interval; (b) the union of two unbounded intervals; (c) the entire real line; (d) an empty set.<sup>8</sup> Case (a) corresponds to a situation where  $\gamma$  is well identified, while (b) and (c) correspond to unbounded confidence sets and indicate (partial or complete) non-identification.

The possibility of getting an empty confidence set may appear surprising. But, on hindsight, this is quite natural: it means that no value of  $\gamma_0$  allows  $\mathscr{H}(\gamma_0)$  to be acceptable. Since  $\mathscr{H}_B$  states there exists a real scalar  $\gamma$  such that  $a_i = (1 - \beta_i)\gamma$ ,  $i = 1, \ldots, n$ , this can be interpreted as a rejection of  $\mathscr{H}_B$ . Further, under  $\mathscr{H}_B$ , the probability that  $CF_{\gamma}(\alpha)$  covers the true value  $\gamma$  is  $1 - \alpha$ , and an empty set obviously does not cover  $\gamma$ . Consequently, the probability that  $CF_{\gamma}(\alpha)$  be empty  $[CF_{\gamma}(\alpha) = \emptyset]$  cannot be greater than  $\alpha$  under  $\mathscr{H}_B : P[CF_{\gamma}(\alpha) = \emptyset] \le \alpha$ . The event  $CF_{\gamma}(\alpha) = \emptyset$  is an exact test with level  $\alpha$  for  $\mathscr{H}_B$  under normality (although its size may be smaller than  $\alpha$ ).

## 4.2. $\mathcal{J}(\gamma_0)$ -based confidence sets

The quadratic confidence set described above relies heavily on the fact that the same critical point  $F_{\alpha}$  can be used to test all values of  $\gamma_0$ . This occurs under the i.i.d. Gaussian distributional assumption, but not necessarily otherwise. Although the quadratic confidence set will remain "asymptotically valid" as long as  $\mathscr{F}(\gamma_0)$  converges to a  $\chi^2(n)$  distribution, this cannot provide an exact confidence set. The Fieller-Hotelling-HAC-type procedure can be extended to allow for possibly non-Gaussian disturbances, by inverting an  $\alpha$ -level test based on  $\mathscr{F}(\gamma_0)$  [or equivalently on  $LR(\gamma_0)$ ] performed by simulation (as a MC test). Consider the MC p-value  $\hat{p}_N(LR;\gamma_0,\nu)$  function associated with this statistic, as defined in (3.11). Since the critical region  $\hat{p}_N(LR;\gamma_0,\nu_0) \leq \alpha$  has size  $\alpha$  for testing  $\gamma = \gamma_0$  when  $\nu_0$  is known, the set of  $\gamma_0$  values for which  $\hat{p}_N(\gamma_0,\nu_0)$  exceeds  $\alpha$ , i.e.

$$C_{\nu}^{LR}(\alpha;\nu) = \left\{ \gamma_0 \in \Gamma : \hat{p}_N(LR;\gamma_0,\nu_0) > \alpha \right\},\tag{4.8}$$

is a confidence set with size  $1-\alpha$  for  $\gamma$ . Similarly, when  $\nu$  is not specified, the test  $\hat{p}_N(LR; \gamma_0, \Omega_{\mathscr{D}}) \leq \alpha$  yields:

$$C_{\mathcal{V}}^{LR}(\alpha; \mathcal{D}) = \left\{ \gamma_0 \in \Gamma : \hat{p}_N(LR; \gamma_0, \Omega_{\mathcal{D}}) > \alpha \right\}, \tag{4.9}$$

whose level is  $1-\alpha$ .  $C_{\gamma}(\alpha; \nu)$  or  $C_{\gamma}(\alpha; \mathcal{D})$  must be drawn by numerical methods. Our empirical analysis reported below, relies on nested grid searches, over  $\gamma_0$  and  $\kappa$ , for the Student-t case (2.15), and over  $\gamma_0$  and  $(\pi, \omega)$  for the normal-mixture case (2.16); for the GARCH case (2.17), we conduct a grid search on  $\gamma_0$  where for each candidate value, we run the simulated annealing optimization algorithm to calculate the maximal p-value from (4.9) over the 2n nuisance parameters in  $\phi$ .

We have no closed-form description of the structure of  $C_{\gamma}^{LR}(\alpha; \nu)$  or  $C_{\gamma}^{LR}(\alpha; \mathcal{D})$ . While these can be bounded intervals (this is shown numerically in Section 8),  $C_{\gamma}^{LR}(\alpha; \nu)$  or  $C_{\gamma}^{LR}(\alpha; \mathcal{D})$  must be unbounded with a high probability if  $\gamma$  is not identifiable or weakly identified (see Dufour, 1997). An empty confidence set is also possible and provides evidence that  $\mathcal{H}_{B}$  is not compatible

<sup>8.</sup> For further discussion, see Dufour and Jasiak (2001); Zivot *et al.* (1998); Dufour and Taamouti (2005); Kleibergen (2009); and Mikusheva (2009).

with the data. The event  $C_{\gamma}^{LR}(\alpha; \nu) = \emptyset$  [or  $C_{\gamma}^{LR}(\alpha; \mathcal{D}) = \emptyset$ ] is a test with level  $\alpha$  for  $\mathcal{H}_B$  under (2.13). When the confidence set  $C_{\gamma}^{LR}(\alpha; \nu)$  is not empty, the identity  $LR(\hat{\gamma}) = \inf \{LR(\gamma_0) : \gamma_0 \in \Gamma\}$  entails that  $\hat{\gamma}$  must belong to  $C_{\gamma}^{LR}(\alpha; \nu)$ .

The Hotelling-based confidence set we obtain for the GARCH case is exact, because the cut-off point we use when inverting  $\mathcal{F}(\gamma_0)$  is adjusted for the parametric form (2.17) via the maximized p-value from (4.9). Inverting  $\mathcal{F}(\gamma_0)$  in (2.28) may however be more appropriate. We thus define using (3.15) and (3.16)

$$C_{\gamma}^{\mathcal{J}}(\alpha;\nu_0) = \left\{ \gamma_0 \in \Gamma : \hat{p}_N(\mathcal{J};\gamma_0,\nu_0) > \alpha \right\},\tag{4.10}$$

which gives a robust confidence set with size  $1-\alpha$  for  $\gamma$ . Similarly, when  $\nu$  is not specified, the test  $\hat{p}_N(LR; \gamma_0, \Omega_{\varnothing}) \le \alpha$  yields

$$C_{\gamma}^{\mathcal{J}}(\alpha;\mathcal{D}) = \left\{ \gamma_0 \in \Gamma : \hat{p}_N(\mathcal{J}; \gamma_0, \Omega_{\mathcal{D}}) > \alpha \right\}, \tag{4.11}$$

a robust confidence set with level  $1-\alpha$ . Again, this must be implemented by numerical methods. Inverting an asymptotic test based on  $\mathcal{J}(\gamma_0)$  can also be considered, relaxing the GARCH restriction. For example, a grid search can be conducted on  $\gamma_0$  where for each candidate value,  $\mathcal{J}(\gamma_0)$  is referred to the  $\chi^2(n)$  distribution; this would circumvent the identification problem asymptotically [as argued e.g. in Stock and Wright (2000)], yet in finite samples, the  $\chi^2(n)$  approximation may perform poorly. Indeed, our simulation results reported below illustrate the severity of this problem. Consequently, we use the MMC method for each candidate  $\gamma_0$ : we maximize over the model parameters as well as over  $\phi$ .

# 5. INVARIANCE AND EXACT DISTRIBUTION OF LRB

In this section, we study the exact distribution of the statistic  $LR_B$ , under both the null hypothesis and the corresponding unrestricted multivariate regression alternative model. Our analysis also provides further information (beyond Theorem 3.1) on distribution of  $LR(\gamma_0)$ . We track and control for the joint role *betas* and scale parameters play in identifying  $\gamma$ .

**Lemma 5.1.** Multivariate scale invariance. The likelihood-ratio statistics  $LR(\gamma_0)$  and  $LR_B$  defined in (2.24) and (2.20) are invariant to replacing  $\tilde{Y}$  by  $\tilde{Y}_* = \tilde{Y}A$ , where A is an arbitrary nonsingular  $n \times n$  matrix.

Such transformations can be viewed as the following affine transformations on Y:

$$Y_* = YA + \tilde{R}_{\mathbf{M}} \iota_n'(I_n - A). \tag{5.1}$$

**Theorem 5.2.** EXACT DISTRIBUTION OF BCAPM LR TESTS. Under (2.1) and (2.13), the distributions of  $LR(\gamma_0)$  and  $LR_B$  depend on (B, K) only through  $\bar{B} = (B - \Delta)K^{-1}$ , and

$$LR(\gamma_0) = T \ln(|\hat{W}(\gamma_0)'\hat{W}(\gamma_0)|/|\hat{W}'\hat{W}|), \quad LR_B = \inf\{LR(\gamma_0): \gamma_0 \in \Gamma\},$$
 (5.2)

where  $\Delta = [0, \iota_n]'$ ,  $\hat{W} = M(X)W$ ,  $\bar{M}(\gamma_0)$  is defined as in (3.1) and

$$\hat{W}(\gamma_0) = \bar{M}(\gamma_0)(X\bar{B} + W) = \bar{M}(\gamma_0)\{\iota_T[a + \gamma_0(\beta - \iota_n)]'K^{-1} + W\}. \tag{5.3}$$

If, furthermore, the null hypothesis  $\mathcal{H}_B$  holds, then

$$\hat{W}(\gamma_0) = (\gamma_0 - \gamma)\bar{M}(\gamma_0)\iota_T(\beta - \iota_n)'K^{-1} + \bar{M}(\gamma_0)W$$
(5.4)

and the distribution of LR<sub>B</sub> depends on (B, K) only through  $\gamma$  and  $(\beta - \iota_n)'K^{-1}$ ; in the Gaussian case (2.14), this distribution involves only one nuisance parameter.

Even though B and K may involve up to  $2n+n^2$  different nuisance parameters [or 2n+n(n+1)/2 parameters, if K is triangular], the latter theorem shows that the number of free parameters in the distributions of  $LR(\gamma_0)$  and  $LR_B$  does not exceed 2n; when  $\mathcal{H}_B$  holds, the number of free parameters is at most n+1. Further, under  $\mathcal{H}(\gamma_0)$  [using (5.4)]  $\bar{B}$  is evacuated, entailing Theorem 3.1. Theorem 5.2 also provides the power function.

### 6. EXACT BOUND PROCEDURES FOR TESTING $\mathscr{H}_{B}$

In this section, we propose tests for  $\mathcal{H}_B$  in the presence of nuisance parameters induced by non-linearity and non-Gaussian error distributions. Because our rationale leads to Shanken, 1986 bound for the Gaussian case, we study first global bounds based on tests of  $\mathcal{H}(\gamma_0)$ . Second, we describe more general but computationally more expensive methods based on the technique of MMC tests to obtain tighter bounds.

# 6.1. Global bound induced by tests of $\mathcal{H}(\gamma_0)$

The results of Section 3 on testing  $\gamma = \gamma_0$  can be used to derive a global bound on the distribution of the statistic  $LR_B$ . This is done in the following theorem.

**Theorem 6.1.** Global bound on the null distribution of the BCAPM test. *Under the assumptions* (2.1), (2.13), and  $\mathcal{H}_B$ , we have, for any given  $v \in \Omega_{\mathscr{D}}$ ,

$$P[LR_{B} \ge x] \le \sup_{\gamma_{0} \in \Gamma} P[\overline{LR}(\gamma_{0}, W) \ge x], \forall x, \tag{6.1}$$

where  $\overline{LR}(\gamma_0, W)$  is defined in (3.1). Further, in the Gaussian case (2.14), we have:

$$P[(T-1-n)(\Lambda_B-1)/n \ge x] \le P[F(n,T-1-n) \ge x], \forall x.$$
 (6.2)

To relate this result to available bounds, observe that (6.1) and (6.2) easily extend to the following multi-beta setups: for i = 1, ..., n, t = 1, ..., T,

$$R_{it} = a_i + \sum_{j=1}^{s} \beta_{ij} \tilde{R}_{jt} + u_{it}, \qquad \mathcal{H}_{B} : a_i = \gamma \left( 1 - \sum_{j=1}^{s} \beta_{ij} \right),$$
 (6.3)

where  $\tilde{R}_{jt}$ ,  $j=1,\ldots,s$ , are returns on s benchmarks. In this case, the bounding distribution of  $LR_B$  obtains as in Theorem 6.1 where  $X=[\iota_T, \tilde{R}_1, \ldots, \tilde{R}_s], \tilde{R}_j=(\tilde{R}_{j1}, \ldots, \tilde{R}_{jT})', j=1,\ldots,s$ , and H is the k-dimensional row vector  $(1, \gamma_0, \ldots, \gamma_0)$ . In the Gaussian case,  $P[\overline{LR}(\gamma_0, W) \ge x]$  does not depend on  $\gamma_0$ , and the bounding distribution under normality is F(n, T-s-n). Shanken (1986) suggests the statistic

$$\hat{Q} = \min_{\gamma} \left\{ \frac{T \left[ \hat{a} - \gamma (\iota_n - \hat{\beta} \iota_s) \right]' \left[ \left( T / (T - 2) \right) \hat{\Sigma} \right]^{-1} \left[ \hat{a} - \gamma (\iota_n - \hat{\beta} \iota_s) \right]}{1 + (\bar{R}_{\mathbf{M}} - \gamma \iota_s)' \hat{\Delta}_{\mathbf{M}}^{-1} (\bar{R}_{\mathbf{M}} - \gamma \iota_s)} \right\}$$
(6.4)

where  $\hat{a}$  is an *n*-dimensional vector that includes the (unconstrained) intercept estimates,  $\hat{\beta}$  is an  $n \times s$  matrix whose rows include the unconstrained OLS estimates of  $(\beta_{i1}, \dots, \beta_{is})$ ,  $i = 1, \dots, n$ ,  $\bar{R}_{M}$  and  $\hat{\Delta}_{M}$  include respectively the time-series means and sample covariance matrix corresponding to the right-hand-side total portfolio returns. Further, the minimum in (6.4) occurs at the constrained MLE  $\hat{\gamma}$  of  $\gamma$ , and

$$LR_{\rm B} = T \ln(1 + \hat{Q}/(T - s - 1)).$$
 (6.5)

For normal errors,  $(T-s-n)\hat{Q}/[n(T-s-1)]$  can be bounded by the F(n, T-n-s) distribution. The latter obtains from Gibbons *et al.*, 1989 joint test of zero intercepts, where returns are expressed in excess of a known  $\gamma$ .

Independently, Stewart (1997) shows [using Dufour (1989)] that, under normal errors,  $(T-s-n)[(|\hat{\Sigma}_{\rm B}|/|\hat{\Sigma}|)-1]/n$  can be bounded by the F(n,T-n-s) distribution. Now, from (2.24) and (6.5), we see that Shanken and Stewart's bounds are equivalent, and both results obtain from Theorem 6.1 in the special case of normal errors.

When disturbances are non-Gaussian, Theorem 6.1 entails that the bounding distribution can easily be simulated. Using the notation we introduced in Section 3.2 [refer in particular to the definitions of the p-value function  $p_N[.]$  from (3.5) and the maximized p-values (3.8)–(3.10)] the following algorithm may be used.

- B1 Given a value of  $\nu$ , generate N i.i.d. draws from the distribution of W then, for any given  $\gamma_0$ , apply the (bounding) function  $\overline{LR}(\gamma_0, W)$  from (3.1) to each draw. With reference to Section 3.2, this implies using  $\overline{LR}(\gamma_0, W)$  for  $\overline{S}(\eta, W)$ , which will yield a vector  $\overline{LR}_N(\gamma_0, \nu)$  of N simulated values of the bounding test statistic.
- B2 Compute the MC p-value [denoted the bound MC (BMC) p-value]

$$\hat{p}_{N}^{\mathrm{U}}(LR_{\mathrm{B}};\gamma_{0},\nu) \equiv p_{N}[LR_{\mathrm{B}}^{(0)}|\overline{LR}_{N}(\gamma_{0},\nu)]. \tag{6.6}$$

In contrast with the Gaussian case,  $\hat{p}_N^U(LR_B; \gamma_0, \nu)$  may depend on  $\gamma_0$ ; nevertheless, for any  $\gamma_0$ ,

$$LR_{\rm B} \leq LR(\gamma_0) \Rightarrow \hat{p}_N(LR; \gamma_0, \nu) \leq \hat{p}_N^{\rm U}(LR_{\rm B}; \gamma_0, \nu). \tag{6.7}$$

B3 Critical regions that provably satisfy the level constraint can be obtained by maximizing  $\hat{p}_N^{\rm U}(LR_{\rm B}; \gamma_0, \nu)$  over the relevant nuisance parameters, leading to the maximized p-values  $\hat{p}_N^{\rm U}(LR_{\rm B}; \Gamma, \nu), \hat{p}_N^{\rm U}(LR_{\rm B}; \gamma_0, \Omega_{\mathscr{D}})$ , and  $\hat{p}_N^{\rm U}(LR_{\rm B}; \Gamma, \Omega_{\mathscr{D}})$ .

**Theorem 6.2.** Global simulation-based bound on the null distribution of the BCAPM TEST STATISTIC. *Under* (2.1), (2.13), and  $\mathcal{H}_{B}$ , we have, using the notation in (3.8)–(3.10):

$$P[\hat{p}_{N}^{U}(LR_{B};\Gamma,\nu) \leq \alpha] \leq \alpha, \quad P[\hat{p}_{N}^{U}(LR_{B};\Gamma,\Omega_{\mathscr{D}}) \leq \alpha] \leq \alpha, \tag{6.8}$$

where v represents the true distributional shape of W.

These bound tests are closely related to the confidence set-based test proposed in Section 4: the null hypothesis is rejected when the confidence set for  $\gamma$  is empty, *i.e.* if no value of  $\gamma_0$  can be deemed acceptable (at level  $\alpha$ ), either with  $\nu$  specified or  $\nu$  taken as a nuisance parameter. This may be seen on comparing (4.9) with the probabilities in Theorem 6.2. Since  $LR_B = \inf\{LR(\gamma_0): \gamma_0 \in \Gamma\}$ , this suggests a relatively easy way of showing that  $C_{\gamma}^{LR}(\alpha; \nu)$  or  $C_{\gamma}^{LR}(\alpha; \mathcal{D})$  is not empty, through the specific p-value  $\hat{p}_N^U(LR_B; \hat{\gamma}, \nu)$  obtained by taking  $\gamma_0 = \hat{\gamma}$  in (6.6). We shall call  $\hat{p}_N^U(LR_B; \hat{\gamma}, \nu)$  the QML-BMC p-value.

**Theorem 6.3.** Relation between efficiency tests and zero-beta confidence sets. *Under* (2.1), (2.13), and  $\mathcal{H}_B$ , let  $\hat{\gamma}$  be the QML estimator of  $\gamma$  in (2.25). Then, using the notation from (3.8) to (3.10):

$$\begin{split} \hat{p}_{N}^{\mathrm{U}}(LR_{\mathrm{B}};\hat{\gamma},\nu) > \alpha \Rightarrow \hat{p}_{N}(LR_{\mathrm{B}};\Gamma,\nu) > \alpha \Rightarrow C_{\gamma}^{LR}(\alpha;\nu) \neq \emptyset, \, \forall \nu \in \Omega_{\mathcal{D}}, \\ \hat{p}_{N}^{\mathrm{U}}(LR_{\mathrm{B}};\hat{\gamma},\Omega_{\mathcal{D}}) > \alpha \Rightarrow \hat{p}_{N}(LR_{\mathrm{B}};\Gamma_{0},\Omega_{\mathcal{D}0}) > \alpha \Rightarrow C_{\gamma}^{LR}(\alpha;\mathcal{D}) \neq \emptyset, \end{split}$$

where  $C_{\nu}^{LR}(\alpha, \nu)$  and  $C_{\nu}^{LR}(\alpha; \mathcal{D})$  are the sets defined in (4.8) and (4.9).

For the Gaussian case, Zhou (1991) and Velu and Zhou (1999) proposed a potentially tighter bound applicable to statistics that can be written as ratios of independent Wishart variables and does not seem to extend easily to other classes of distributions. In the next section, we propose an approach that yields similarly tighter bounds for non-Gaussian distributions as well. Finally, the HAC statistic  $\mathcal{J}_B$  may be used to obtain alternative identification-robust bound tests following the same rationale. The correspondence between such tests and empty confidence sets entailed by test inversion also follows from similar arguments. Finite-sample MMC level corrections are recommended, given the simulation results in Section 7.

## 6.2. Tighter bounds

Another approach to testing  $\mathcal{H}_B$  with the statistic  $LR_B$  consists in directly assessing its dependence on nuisance parameters and adjusting the test accordingly through the MMC method (Dufour, 2006). Let  $\theta = \psi(B, K)$  represent the parameter vector upon which the distribution of  $LR_B$  actually depends, and  $\Omega_B$  the set of admissible values for  $\theta$  under  $\mathcal{H}_B$ . The dimension of  $\theta$  may be lower than than the number of parameters in B and K. Define the function  $\overline{LR}_B(\theta, W) = \overline{LR}_B(\psi(B,K),W)$  that assigns to each value of (B,K) and the noise matrix W the following outcome: using  $\theta$  and a draw from the distribution of W (which may depend on  $\nu$ ), generate a sample from (2.1) to (2.2), and compute  $LR_B$  [as defined in (2.24)] from this sample. This suggests we can use  $\overline{LR}_B(\theta,W)$  for  $\overline{S}(\eta,W)$ . Maintaining further notation from Section 3.2 [again, recall in particular the definitions of the p-value function  $p_N[.]$  from (3.5) and the maximized p-values (3.8)–(3.10)] the following algorithm may be used.

- C1 Generate N independent replications of W and apply  $\overline{LR}_{B}(\theta, W)$  to each draw, for any value of  $\theta$ , leading to a vector  $\overline{LR}_{BN}(\theta, \nu)$  of N simulated statistics.
- C2 Compute the MC *p*-value

$$\hat{p}_{N}^{\mathrm{B}}(LR_{\mathrm{B}};\theta,\nu) = p_{N} \left[ LR_{\mathrm{B}}^{(0)} \middle| \overline{LR}_{\mathrm{B}N}(\theta,\nu) \right].$$

C3 Critical regions that provably satisfy the level constraint can be obtained by maximizing  $\hat{p}_N^B(LR_B;\theta,\nu)$  over relevant nuisance parameters leading to  $\hat{p}_N^B(LR_B;\Omega_B,\nu)$ ,  $\hat{p}_N^B(LR_B;\theta,\Omega_{\mathscr{D}})$ , and  $\hat{p}_N^B(LR_B;\Omega_B,\Omega_{\mathscr{D}})$ .

Theorem 6.3 guarantees that  $\hat{p}_N^U(LR_B;\Gamma,\nu) \leq \alpha \Rightarrow \hat{p}_N^B(LR_B;\Omega_B,\nu) \leq \alpha$  for any given  $\nu$ . So it may be useful to check the global bound for significance before turning to the MMC one. Furthermore, it is not always necessary to run the numerical maximization underlying MMC to convergence: if  $\hat{p}_N^B(LR_B;\theta,\nu) > \alpha$  given any relevant  $\theta$  (or  $\nu$ ), then a non-rejection is confirmed. We suggest to use the QML estimate  $\hat{\theta}$  of  $\theta$  as start-up value, because this provides *parametric* 

bootstrap-type [or a local MC (LMC)] p-values:

$$p_N^b(LR_B; \nu) = \hat{p}_N^B(LR_B; \hat{\theta}, \nu), \quad p_N^b(LR_B; \Omega_{\mathscr{D}}) = \hat{p}_N^B(LR_B; \hat{\theta}, \Omega_{\mathscr{D}}). \tag{6.9}$$

Then  $p_N^b(LR_B; \nu) > \alpha$  entails  $\hat{p}_N^B(LR_B; \Omega_B, \nu) > \alpha$ , and  $p_N^b(LR_B; \Omega_{\mathscr{D}}) > \alpha$  entails  $\hat{p}_N^B(LR_B; \Omega_B, \Omega_{\mathscr{D}}) > \alpha$ .

Following the same reasoning, a parametric MMC test imposing (2.17) may be applied to the HAC statistics  $\mathcal{J}(\gamma_0)$  and  $\mathcal{J}_B$ , as an attempt to correct their size for the GARCH alternative of interest. We investigate the size-corrected power associated with these statistics in Section 7.

Weak identification of  $\gamma$  implies that the data is weakly informative on the parameter and thus on the related specification. This means that the data may not provide sufficient information on whether  $\mathcal{H}_B$  should be refuted or not. When  $\gamma$  is weakly identified, tests for  $\mathcal{H}_B$  will suggest accepting the underlying pricing restrictions. We thus recommend to interpret tests for  $\mathcal{H}_B$  in conjunction with associated confidence sets that will be unbounded when  $\gamma$  is weakly identified and will provide a much more complete statistical analysis.

## 6.3. Two-stage bound confidence procedures

To deal with the fact that the distribution of W may involve an unknown parameter  $v \in \Omega_{\mathscr{D}}$ , we suggested above to maximize the relevant p-values over  $\Omega_{\mathscr{D}}$ . We next consider restricting the maximization over v to a set that is empirically relevant, as in Beaulieu et al. (2007). This leads to two basic steps: (i) an exact confidence set with level  $1-\alpha_1$  is built for v, and (ii) the MC p-values (presented above) are maximized over all values of v in the latter confidence set and are referred to level  $\alpha_2$ , so that the global test level is  $\alpha = \alpha_1 + \alpha_2$ . In our empirical application, we used  $\alpha/2$ . Let  $\mathscr{C}_v(\alpha_1) = \mathscr{C}_v(\alpha_1; Y)$  be a confidence set with level  $1-\alpha_1$  for v. Then, under  $\mathscr{H}(\gamma_0)$ , we have  $P[\hat{p}_V^V(LR_B; \gamma_0, \mathscr{C}_v(\alpha_1)) \leq \alpha_2] \leq \alpha_1 + \alpha_2$  while, under  $\mathscr{H}_B$ :

$$P[\hat{p}_{N}^{U}(LR_{B};\Gamma,\mathscr{C}_{\nu}(\alpha_{1})) \leq \alpha_{2}] \leq \alpha_{1} + \alpha_{2}, \qquad P[\hat{p}_{N}^{B}(LR_{B};\Omega_{B},\mathscr{C}_{\nu}(\alpha_{1})) \leq \alpha_{2}] \leq \alpha_{1} + \alpha_{2}. \quad (6.10)$$

Note also that for  $\hat{p}_N^B(LR_B; \Omega_B, \mathcal{C}_{\nu}(\alpha_1)) \leq \alpha_2$  not to hold, the following condition is sufficient:

$$\hat{p}_{N}^{B}(LR_{B}; \hat{\theta}, \mathcal{C}_{\nu}(\alpha_{1})) > \alpha_{2}. \tag{6.11}$$

To build a confidence sets for  $\nu$ , we invert a test (of level  $\alpha_1$ ) for the specification underlying (2.13) where  $\nu = \nu_0$  for known  $\nu_0$ ; this avoids the need to use regularity assumptions on  $\nu$ . The test we invert is the three-stage MC goodness-of-fit test introduced in Dufour *et al.* (2003), which uses Mardia, 1970 well-known multivariate skewness and kurtosis measures (SK and KU below):

$$CSK(\nu_0) = 1 - \min \{ \hat{p}_N(ESK(\nu_0); \nu_0), \hat{p}_N(EKU(\nu_0); \nu_0) \}$$
(6.12)

$$ESK(\nu_0) = |SK - \overline{SK}(\nu_0)|, \quad EKU(\nu_0) = |KU - \overline{KU}(\nu_0)|$$
(6.13)

$$SK = \frac{1}{T^2} \sum_{t=1}^{T} \sum_{i=1}^{T} \hat{d}_{ii}^3, \quad KU = \frac{1}{T} \sum_{t=1}^{T} \hat{d}_{tt}^2,$$
 (6.14)

where  $\hat{d}_{it}$  are the elements of the matrix  $\hat{U}(\hat{U}'\hat{U}/T)^{-1}\hat{U}'$ ,  $\overline{SK}(\nu_0)$ , and  $\overline{KU}(\nu_0)$  are simulation-based estimates of the expected SK and KU given (2.13) and  $\hat{p}[ESK(\nu_0)]$  and  $\hat{p}[EKU(\nu_0)]$  are p-values, obtained by MC methods under (2.13). The test's three stages (summarized below) are motivated by the following [see Dufour *et al.* (2003) for proofs and more detailed algorithms].

Let  $\bar{d}_{it}(\nu)$  refer to the elements of the matrix  $\hat{W}(\hat{W}'\hat{W}/T)^{-1}\hat{W}'$  with  $\hat{W}=M(X)W$  and recall that  $\hat{U}=M(X)U=M(X)WK=\hat{W}K$ . Then under assumption (2.13) [including in particular the distributions from (2.14) to (2.16)], SK and KU depend on the data only via  $\bar{d}_{it}(\nu)$  and thus may easily be simulated given draws from the distribution of W. The three stages of the test use this result, as follows.

- D1 Stage I, to obtain  $\overline{SK}(\nu_0)$  and  $\overline{KU}(\nu_0)$ : for a given  $\nu_0$ , draw  $N_0$  samples from the hypothesized distribution of W. Applying  $\bar{d}_{it}(.)$  to each draw, and on replacing  $\hat{d}_{it}$  in (6.14) by  $\bar{d}_{it}(\nu)$  compute the corresponding measures of skewness and kurtosis, and then take their average. The remaining two stages condition on these estimates. Refer to  $\overline{SK}(\nu_0)$  and  $\overline{KU}(\nu_0)$  so obtained as the *reference simulated moments* [**RSM**].
- D2 Stage II, to derive  $CSK(\nu_0)$ : first obtain  $\hat{p}_N(ESK(\nu_0); \nu_0)$  and  $\hat{p}_N(EKU(\nu_0); \nu_0)$ , using the observed statistics and  $N_1$  [where  $\alpha(N_1+1)$  is an integer] draws from the hypothesized distribution of W (independently from stage I), as in the algorithms A1–A3 from Section 3.2 and using the same RSM for the observed and the simulated values of SK and KU. Refer to the N simulated SK and KU from step A1 as the reference simulated series (**RSS**). Applying (6.12) gives the observed value of the test statistic,  $CSK(\nu_0)^{(0)}$ .
- D3. Stage III, to derive a MC p-values for  $CSK(v_0)$ : independently of the previous RSM and BSS, generate N [where  $\alpha(N+1)$  is an integer] additional i.i.d. realizations of W. Applying  $\bar{d}_{it}(.)$  to each draw, and replacing  $\hat{d}_{it}$  in (6.14) by  $\bar{d}_{it}(v)$  compute the corresponding measures of excess skewness and kurtosis (using the RSM). By referring each of the latter to the RSS, compute MC p-values: in other words, repeat step D2 replacing the observed statistics by those last simulated. These p-values provide a vector  $\overline{CSK}_N(v_0)$  of N replications of the combined statistic, leading in turn to  $\hat{p}_N(CSK(v_0); v_0) = p_N[CSK(v_0)^{(0)}|\overline{CSK}_N(v_0)]$ .

The confidence set for  $\nu$  corresponds to the values of  $\nu_0$  that are not rejected at level  $\alpha_1$ , using the latter p-value. For the GARCH case, pre-estimating the  $2n \times 1$  vector  $\phi$  is infeasible with 5 or even 10-year subsamples of monthly data. Nevertheless, the single stage MMC is valid despite this limitation. Interestingly, the simulation study we report next suggests that power costs are unimportant even with relatively small samples.

## 7. SIMULATION STUDY

We now present a small simulation study to assess the performance of the proposed methods. The design is calibrated to match our empirical analysis (see Section 8) that relies on monthly returns of 12 portfolios of NYSE firms over 1927–1995. We consider model (2.1) where  $\tilde{R}_{Mt}$ ,  $t=1,\ldots,T$ , are the returns on the market portfolio from the aforementioned data over the last 5-and 10-year subperiods, as well as the whole sample. We thus take n=12 and T=60,120, and 828. The coefficients of (2.1) including  $\gamma$  are set to their QML estimates (restricted under  $\mathcal{H}_B$  over the conformable sample period). From the QML regression, we also retain the estimated error covariance matrix, to generate model shocks; formally, we compute the corresponding empirical Cholesky factor (denoted  $\hat{K}$ ) and use it for K in (2.13). Test sizes with  $K=I_{12}$  are also analysed to illustrate the effects of portfolio repacking.

We consider normal and Student t-errors (with  $\kappa = 8$ , in accordance with the kurtosis observed in the empirical application), so the random vectors  $W_t$ , t = 1, ..., T, in (2.13) are generated

following (2.14) and (2.15), respectively. The MC tests are applied imposing and ignoring information on  $\kappa$ , which allows us to document the cost of estimating this parameter. When  $\kappa$  is considered unknown, MMC p-values are calculated over the interval  $4 \le \kappa \le 13$  to keep execution time manageable (a wider range is allowed for the empirical application in Section 8). We also consider the case of GARCH errors (2.17), with  $\phi_{1i} = \phi_1$  and  $\phi_{2i} = \phi_2$ ,  $i = 1, \ldots, n$  (the coefficients are the same across equations). This restriction is motivated by execution time, but it is relaxed in Section 8. We use the diagonal elements of  $\hat{K}\hat{K}'$  to scale the intercept, yet we also consider the case where  $\sigma_i^2 = 1$ ,  $i = 1, \ldots, n$ . Samples are simulated with  $(\phi_1, \phi_2) = (.15, .80)$ . These parameters are treated, in turn, as known and as unknown quantities. In view of the low dimension of the nuisance parameter space in this case, when  $(\phi_1, \phi_2)$  is treated as unknown, p-value maximization is achieved through a coarse grid search (for the purpose of this simulation). The p-value function does not appear to be very sensitive to the value of  $(\phi_1, \phi_2)$ , and the results presented below indicate this is sufficient for controlling test level in the relevant cases. A more thorough optimization is however used in Section 8.

The results of the simulation are summarized in Tables 1–3. These tables report empirical rejection rates for various tests of  $\mathcal{H}(\gamma_0)$  with nominal size 5%. These rejection rates determine the coverage properties of confidence sets derived from the tests. Since we focus on estimating  $\gamma$ ,  $\mathcal{H}_B$  is imposed for both the size and power studies. We compare the following tests: (1) a Wald-type test that rejects  $\gamma = \gamma_0$  when  $\gamma_0$  falls outside the Wald-type confidence interval  $[\hat{\gamma} - 1.96 \times \text{AsySE}(\hat{\gamma}), \hat{\gamma} + 1.96 \times \text{AsySE}(\hat{\gamma})]$ , using the QML estimator  $\hat{\gamma}$ , an asymptotic standard

TABLE 1
Tests on zero-beta rate: empirical size

n=12			T =	=60	T =	120	T =	828
				K		K		K
W	Statistic	<i>p</i> -value	$I_{12}$	ĥ	$I_{12}$	ĥ	$I_{12}$	ĥ
Normal	Wald-type	N[0,1]	0.709	0.196	0.633	0.096	0.578	0.050
	$LR(\gamma_0)$	MC	0.057	0.057	0.048	0.048	0.041	0.041
Student-t	Wald-type	N[0,1]	0.714	0.218	0.645	0.106	0.587	0.055
	$LR(\gamma_0)$	MC: $\kappa$ known	0.053	0.053	0.046	0.046	0.043	0.043
		MMC: $\kappa$ unknown	0.043	0.043	0.035	0.035	0.031	0.031
GARCH	Wald-type	N[0,1]	0.676	0.200	0.628	0.086	0.579	0.047
	$LR(\gamma_0)$	MC: $\phi_1 = \phi_2 = 0$	0.059	0.059	0.048	0.048	0.046	0.046
		MC: $\phi_1, \phi_2$ known	0.064	0.064	0.043	0.043	0.050	0.028
		MMC: $\phi_1, \phi_2$ unknown	0.054	0.054	0.032	0.032	0.028	0.050
	$\mathscr{J}(\gamma_0)$	$\chi^{2}(12)$	0.954	0.954	0.686	0.686	0.127	0.127
	V	MC: $\phi_1, \phi_2$ known	0.049	0.049	0.045	0.045	0.049	0.040
		MMC: $\phi_1, \phi_2$ unknown	0.040	0.040	0.034	0.034	0.040	0.049

Notes: The table reports the empirical rejection rates of various tests for  $\mathcal{H}(\gamma_0)$  with nominal level 5%. The values of  $\gamma_0$  tested are:  $\gamma_0 = -0.000089$  for T = 60,  $\gamma_0 = 0.004960$  for T = 120,  $\gamma_0 = 0.005957$  for T = 828. The design is calibrated to match our empirical analysis (see Section 8). The tests compared are the following. (1) A Wald-type test that rejects  $\gamma = \gamma_0$  when  $\gamma_0$  falls outside the Wald-type confidence interval  $[\hat{\gamma} - 1.96 \times \text{AsySE}(\hat{\gamma}), \hat{\gamma} + 1.96 \times \text{AsySE}(\hat{\gamma})]$ , using the QML estimator  $\hat{\gamma}$  with asymptotic standard error [AsySE( $\hat{\gamma}$ )] based on (2.27), and a normal limiting distribution. (2) MC and MMC tests based on  $LR(\gamma_0)$  in (2.20), with MC p-values for i.i.d. normal and Student-t errors (with known or unknown  $\kappa$ ), Gaussian GARCH with known or unknown ( $\phi_1, \phi_2$ ), as well as  $\phi_1 = \phi_2 = 0$  (i.e. ignoring the GARCH dependence even when it is present in the simulated process). (3) Tests based on the HAC Wald-type statistic  $\mathcal{J}(\gamma_0)$  in (2.28), using a  $\chi^2(n)$  critical value, MC with known ( $\phi_1, \phi_2$ ), and MMC where ( $\phi_1, \phi_2$ ) is taken as unknown. In the i.i.d. cases, the errors are generated using (2.13) with K set to either  $I_{12}$  or  $\hat{K}$ , which corresponds to the Cholesky factor of the least-squares error covariance estimate from the empirical data used for the simulation design. In the GARCH case, samples are generated with conditional variance as in (2.17) using  $\hat{K}$  or  $I_{12}$  for K.

		•	•			_		
n=12		T	=60	T = 120		T = 828		
$\overline{W}$	Statistic	<i>p</i> -value	Step	Power	Step	Power	Step	Power
Normal	$LR(\gamma_0)$	MC: $\phi_1 = \phi_2 = 0$	0.50	0.151	0.50	0.226	0.20	0.129
			0.75	0.315	0.75	0.529	0.30	0.313
			1.0	0.544	1.0	0.835	0.50	0.814
			2.0	0.981	1.5	0.999	0.75	0.998
Student-t	$LR(\gamma_0)$	MC: κ known	0.50	0.134	0.50	0.181	0.20	0.109
		MMC: $\kappa$ unknown		0.126		0.158		0.080
		MC: $\kappa$ known	0.75	0.264	0.75	0.428	0.30	0.237
		MMC: $\kappa$ unknown		0.239		0.384		0.182
		MC: $\kappa$ known	1.0	0.494	1.0	0.709	0.50	0.660
		MMC: $\kappa$ unknown		0.440		0.673		0.605
		MC: $\kappa$ known	2.0	0.939	1.5	0.997	0.75	0.966
		MMC: k unknown		0.925		0.997		0.960

TABLE 2
Tests on zero-beta rate: empirical power. Gaussian and Student designs

Notes: The table reports the empirical rejection rates of various tests for  $\mathcal{H}(\gamma_0)$  with nominal level 5%. The values of  $\gamma_0$  tested are:  $\gamma_0 = -0.000089$  for T = 60,  $\gamma_0 = 0.004960$  for T = 120,  $\gamma_0 = 0.005957$  for T = 828. The sampling design conforms with the size study, for the  $\hat{K}$  case. Samples are drawn with  $\gamma$  calibrated to its QML counterpart from the 1991–1995 subsample; values for  $\gamma_0$  are set to the latter value +  $step \times \hat{\sigma}_i^{min}$ , where  $\hat{\sigma}_i^{min} = [min\{\hat{\sigma}_i^2\}]^{1/2}$ , and  $\hat{\sigma}_i^2$  are the diagonal terms of  $\hat{K}\hat{K}'$ . See Table 1 for further details on the design and tests applied.

TABLE 3
Tests on zero-beta rate: empirical power. Gaussian GARCH design

	n=12	T	=60	T =	= 120	T = 828	
Statistic	<i>p</i> -value	Step	Power	Step	Power	Step	Power
$LR(\gamma_0)$	MC: $\phi_1 = \phi_2 = 0$	0.50	0.112	0.50	0.203	0.20	0.195
	MC: $\phi_1, \phi_2$ known		0.113		0.204		0.198
	MMC: $\phi_1, \phi_2$ unknown		0.106		0.170		0.168
$\mathscr{J}(\gamma_0)$	MMC: $\phi_1, \phi_2$ known		0.088		0.155		0.208
	MMC: $\phi_1, \phi_2$ unknown		0.078		0.133		0.183
$LR(\gamma_0)$	MC: $\phi_1 = \phi_2 = 0$	0.75	0.247	0.75	0.449	0.30	0.465
	MC: $\phi_1, \phi_2$ known		0.248		0.452		0.471
	MMC: $\phi_1, \phi_2$ unknown		0.213		0.411		0.425
$\mathscr{J}(\gamma_0)$	MMC: $\phi_1, \phi_2$ known		0.177		0.316		0.442
	MMC: $\phi_1, \phi_2$ unknown		0.158		0.276		0.411
$LR(\gamma_0)$	MC: $\phi_1 = \phi_2 = 0$	1.0	0.447	1.0	0.753	0.50	0.945
	MC: $\phi_1, \phi_2$ known		0.441		0.753		0.950
	MMC: $\phi_1, \phi_2$ unknown		0.395		0.709		0.937
$\mathscr{J}(\gamma_0)$	MC: $\phi_1, \phi_2$ known		0.300		0.552		0.934
	MMC: $\phi_1, \phi_2$ unknown		0.269		0.505		0.920
$LR(\gamma_0)$	$MC, \phi_1 = \phi_2 = 0$	2.0	0.913	1.5	.973	0.75	1.0
	MC: $\phi_1, \phi_2$ known		0.915		0.970		1.0
	MMC: $\phi_1, \phi_2$ unknown		0.892		0.962		1.0
$\mathscr{J}(\gamma_0)$	MMC: $\phi_1, \phi_2$ known		0.719		0.856		1.0
0	MMC: $\phi_1, \phi_2$ unknown		0.664		0.931		1.0

Notes: The values of  $\gamma_0$  tested are:  $\gamma_0 = -0.000089$  for T = 60,  $\gamma_0 = 0.004960$  for T = 120,  $\gamma_0 = 0.005957$  for T = 828. Numbers reported are empirical rejection rates for various tests of  $\mathcal{H}(\gamma_0)$  with nominal size 5%. The sampling design conforms with the size study, for the  $\hat{K}$  case; errors are generated with conditional variance as in (2.17) using  $\hat{K}$ . See Table 1 for a complete description of the designs and tests applied. Samples are drawn with  $\gamma$  calibrated to its QML counterpart from the 1991–1995 subsample; values for  $\gamma_0$  are set to the latter value +  $step \times \hat{\sigma}_i^{min}$  (for various step values) where  $\hat{\sigma}_i^{min} = [min\{\hat{\sigma}_i^2\}]^{1/2}$ , and  $\hat{\sigma}_i^2$  are the diagonal terms of  $\hat{K}\hat{K}'$ .

error [AsySE( $\hat{\gamma}$ ), based on (2.27)], and a normal limiting distribution; (2) the MC and MMC tests based on the QLR test statistic  $LR(\gamma_0)$  defined in Theorem 3.1, with MC p-values for i.i.d. normal or Student-t errors (with known or unknown  $\kappa$ ), Gaussian GARCH with known or unknown  $(\phi_1, \phi_2)$ , as well as as  $\phi_1 = \phi_2 = 0$  (i.e. ignoring the GARCH dependence even when it is present in the simulated process); (3) tests based on the HAC Wald-type statistic  $\mathcal{J}(\gamma_0)$  in (2.28), using a  $\chi^2(n)$  critical value, MC with known  $(\phi_1, \phi_2)$ , and MMC where  $(\phi_1, \phi_2)$  is taken as unknown.

In the size study (Table 1),  $\gamma_0$  is calibrated to its QML counterpart from the data set  $[\gamma_0 = -0.000089 \, \text{for} \, T = 60, \gamma_0 = 0.004960 \, \text{for} \, T = 120, \gamma_0 = 0.005957 \, \text{for} \, T = 828]$ . In the power study (Tables 2 and 3), we focus on the  $\hat{K}$  case; samples are drawn with  $\gamma$  set to its QML estimate, and  $\gamma_0$  is set to the latter value +  $step \times \widehat{\sigma}_i^{\min}$ , where  $\widehat{\sigma}_i^{\min} = [\min\{\widehat{\sigma}_i^2\}]^{1/2}$ , and  $\widehat{\sigma}_i^2$  are the diagonal terms of  $\hat{K}\hat{K}'$  (with various step values). N = 99 is used for MC tests (N = 999 is used in the empirical application). In each experiment, the number of simulations is 1000. We use 12 lags for the HAC correction.

Our results can be summarized as follows. The asymptotic i.i.d. or robust procedures are very unreliable from the viewpoint of controlling level. Whereas we observe empirical frequencies of type I errors over 70% and sometimes 90% with T=60, we still see empirical rejections near 55% with T=828. The results also show that the empirical size of the HAC-based tests is not affected by K, though a formal proof of its invariance is not available. This observation is however compatible with the fact that its size improves with larger samples: while the level of the Wald-type test shows no improvement (around 55%) even with T=828 and normal errors, the size of the  $\mathcal{J}(\gamma_0)$  statistic drops from 95% with T=60 to 12% with T=828. The likelihood-ratio and robust MC and MMC tests achieve level control; in the GARCH case, the MC likelihood-ratio test has the correct size even when GARCH dependence is not accounted for.

In view of the poor size performance of the asymptotic tests, the power study focuses on procedures whose level appears to be under control. Overall, the MMC correction is not too costly from the power viewpoint, with both Student-t and GARCH errors. In the latter case, the likelihood-ratio-type test uncorrected for GARCH effects outperforms all the other tests. When GARCH corrections are performed via MMC, the likelihood-ratio-type test performs generally better than the  $\mathcal{J}(\gamma_0)$  test.

#### 8. EMPIRICAL ANALYSIS

In this section, we assess  $\mathcal{H}_B$  as defined in (2.2) in the context of (2.1) under the distributional assumptions (2.15) and (2.16), as well as the Gaussian GARCH in (2.17). We use real monthly returns over the period going from January 1926 to December 1995, obtained from CRSP. The data studied involve 12 portfolios of NYSE firms grouped by standard two-digit industrial classification (SIC). The sectors studied include: (1) petroleum; (2) finance and real estate; (3) consumer durables; (4) basic industries; (5) food and tobacco; (6) construction; (7) capital goods; (8) transportation; (9) utilities; (10) textile and trade; (11) services; (12) leisure; for details on the SIC codes, see Beaulieu *et al.* (2007). For each month, the industry portfolios include the firms for which the return, the price per common share and the number of shares outstanding are recorded by CRSP. Furthermore, portfolios are value weighted in each month. We measure the market return by value weighted NYSE returns, and the real risk-free rate by the one-month Treasury bill rate net of inflation, both available from CRSP. All MC tests were applied with N=999 and MMC p-values are obtained using the simulated annealing algorithm.

Our QML-based BCAPM test results are summarized in Table 4. Non-Gaussian p-values are the largest MC p-values over the error distribution parameters [respectively:  $\kappa$  and  $(\pi, \omega)$  for (2.15) and (2.16)] within the specified 97.5% confidence sets; the latter are reported in Table 5. In the GARCH case (2.17), p-values are the largest MC p-values over all  $(\phi_{1i}, \phi_{2i})$ . Given a

TABLE 4
QML-based tests of BCAPM

Sample	$LR_{ m B}$	$p_{\infty}$		Normal		GARCH
			LMC	MMC	BND	BND
1927 – 30	16.10	0.137	0.269	0.308	0.366	0.340
1931 - 35	14.09	0.228	0.344	0.381	0.432	0.451
1936 - 40	15.36	0.167	0.257	0.284	0.345	0.355
1941 - 45	18.62	0.068	0.148	0.163	0.203	0.213
1946 - 50	32.69	0.001	0.005	0.006	0.007	0.006
1951 - 55	37.04	0.000	0.003	0.004	0.004	0.003
1956 - 60	26.10	0.006	0.027	0.031	0.042	0.039
1961 - 65	29.21	0.002	0.011	0.016	0.020	0.015
1966 - 70	27.45	0.004	0.016	0.018	0.026	0.029
1971 - 75	16.81	0.113	0.213	0.224	0.292	0.294
1976 - 80	25.76	0.007	0.027	0.031	0.040	0.042
1981 - 85	14.98	0.183	0.316	0.335	0.387	0.404
1986 - 90	35.41	0.000	0.003	0.004	0.004	0.005
1991-95	16.41	0.127	0.219	0.253	0.310	0.320
		Student-t			Normal mixtur	e
	LMC	MMC	BND	LMC	MMC	BND
1927 – 30	0.272	0.316	0.360	0.279	0.313	0.381
1931 - 35	0.359	0.399	0.468	0.342	0.387	0.452
1936 - 40	0.282	0.308	0.372	0.265	0.302	0.357
1941 - 45	0.147	0.169	0.210	0.150	0.165	0.211
1946 - 50	0.007	0.007	0.010	0.007	0.007	0.008
1951 - 55	0.003	0.005	0.005	0.003	0.003	0.003
1956 - 60	0.030	0.040	0.052	0.028	0.035	0.045
1961 - 65	0.013	0.017	0.023	0.014	0.021	0.024
1966 - 70	0.020	0.025	0.032	0.018	0.023	0.028
1971 - 75	0.217	0.248	0.300	0.206	0.238	0.292
1976 - 80	0.026	0.035	0.039	0.026	0.034	0.042
1981 - 85	0.323	0.399	0.405	0.318	0.339	0.406
1986 - 90	0.004	0.005	0.005	0.004	0.004	0.005
1991 - 95	0.226	0.263	0.325	0.226	0.261	0.319

Notes:  $LR_B$  is the statistic in (2.24). Remaining numbers are associated p-values.  $p_\infty$  is based on  $\chi^2(n-1)$ . All other non-Gaussian p-values are the largest MC p-values over the shape parameter  $\nu$  within the specified confidence sets [ $\nu = \kappa$  or  $\nu = (\pi, \omega)$ ; refer to Table 5]. LMC is the bootstrap p-value in (6.11) and MMC is the maximal p-value in (6.10) (refer to Section 6.2). BND is the bound (6.2) for the Gaussian case and the QML-BMC bound from Theorem 6.3 otherwise; the GARCH BND is the largest QML-BMC over  $\phi_{1i}, \phi_{2i}$  [from (2.17)]. Returns for the months of January and for October 1987 are excluded. Given a 5% level, the cut-off is 0.05 for  $p_\infty$ , the normal LMC, MMC, and the GARCH p-values; for the Student-t and mixtures, the cut-off is 0.025. p-values which lead to significant tests with this benchmark are in bold.

5% level test, the benchmark is .05 for  $p_{\infty}$ , normal LMC, MMC, and GARCH p-values, while the Student-t and normal mixture p-values should be compared to 0.025 (to ensure that the test has level 0.05). Non-rejections by LMC MC p-values are conclusive (though rejections are not); rejections based on the conservative bound reported under the heading BND are conclusive under normality; non-rejections based on the QML bound in the non-Gaussian case (reported under the heading BND) signal that the confidence set for  $\gamma$  is not empty; however, since the MMC p-value is based on the tightest bound, this evidence does not necessarily imply non-rejection of  $\mathcal{H}_{\rm B}$ .

The empirical results presented in Table 4 show that the asymptotic test and the Gaussian-based bound test yield the same decision at level 5%, although the former p-values are much lower. The non-normal p-values exceed the Gaussian-based p-value, enough to change the test decision. For instance, at the 5% significance level, we find seven rejections of the null hypothesis for the

		Mixture $(\pi, \omega)$ , confidence set for $\omega$						
	(1)	(2)	(3)	(4)	(5)	(6)		
Sample	$\pi = 0.1$	$\pi = 0.2$	$\pi = 0.3$	$\pi = 0.4$	$\pi = 0.5$	κ		
1927 – 30	≥ 1.8	1.6-2.8	1.6-2.5	1.6-2.5	1.6-2.6	3-12		
1931 - 35	2.1 - 10.0	1.9 - 3.0	1.9 - 2.7	1.9 - 2.7	2.1 - 3.0	3 - 8		
1966 - 40	1.5 - 3.5	1.5 - 2.3	1.4 - 2.1	1.4 - 2.0	1.4 - 2.1	4 - 25		
1941 - 45	1.3 - 3.5	1.3 - 2.1	1.3 - 1.9	1.3 - 1.8	1.3 - 1.9	≥5		
1946 - 50	1.4 - 3.5	1.3 - 2.2	1.3 - 2.0	1.3 - 1.9	1.3 - 1.9	5 - 37		
1951 - 55	1.4 - 3.5	1.4 - 2.2	1.3 - 2.0	1.3 - 1.9	1.3 - 2.0	5 - 34		
1956 - 60	1.3 - 2.8	1.2 - 2.0	1.2 - 1.9	1.2 - 1.8	1.2 - 1.8	≥5		
1961 - 65	1.0 - 2.2	1.0 - 1.6	1.0 - 1.5	1.0 - 1.5	1.0 - 1.5	_ ≥7		
1966 - 70	1.3 - 3.0	1.3 - 2.0	1.3 - 1.9	1.3 - 1.8	1.2 - 1.9	_ ≥5		
1971 - 75	1.5 - 3.5	1.5 - 2.2	1.4 - 2.0	1.4 - 1.9	1.4 - 2.0	4-24		
1976 - 80	1.6 - 4.0	1.5 - 2.5	1.5 - 2.2	1.5 - 2.2	1.5 - 2.3	4 - 19		
1981 - 85	1.4 - 3.5	1.4 - 2.1	1.3 - 2.0	1.3 - 1.9	1.4 - 2.0	5 - 33		
1986 - 90	1.1 - 3.0	1.1 - 2.0	1.1 - 1.8	1.0 - 1.7	1.1 - 1.8	≥5		
1991 - 95	1.0 - 1.9	1.0 - 1.5	1.0 - 1.4	1.0 - 1.3	1.0 - 1.3	_ ≥19		

TABLE 5
Confidence sets for intervening parameters

Notes: Numbers in Columns (1)–(5) represent a confidence set for the parameters  $(\pi, \omega)$  [respectively, the probability of mixing and the ratio of scales] of the multivariate mixtures-of-normal error distribution. Column (6) presents the confidence set for  $\kappa$ , the degrees-of-freedom parameter of the multivariate Student-t error distribution. See Section 6 for details on the construction of these confidence sets: the values of  $(\pi, \omega)$  or  $\kappa$  (respectively) in this set are not rejected by the CSK test (6.12) (see Dufour et al., 2003) under multivariate mixtures or Student-t errors (respectively). Note that the maximum of the p-value occurs in the closed interval for  $\omega$ . Returns for the month of January and October 1987 are excluded from the data set.

asymptotic  $\chi^2(11)$  test, seven for the MC *p*-values under normality and with normal GARCH, and five (relying on the MMC *p*-value) under the Student-*t* and normal mixture distributions.

Focusing on Student-*t* and normal mixture distributions with parameters not rejected by proper GF tests, we see that mean-variance efficiency test results can change relative to the available *F*-based bound. The power advantages of the MMC procedure are illustrated by the results of the 1966–1970 subperiod where the QML *p*-value exceeds 2.5% for the Student-*t* and normal mixture distributions, whereas the MMC *p*-value signals a rejection.

The confidence sets for distributional parameters are reported in Table 5. In the mixture case, the confidence region is summarized as follows for presentation ease: we give the confidence set for  $\omega$  corresponding to five different values of  $\pi$ .

In Table 6, we present: (i) the average real market return as well as the average real risk-free rate over each subperiod, (ii) the QML estimate of  $\gamma$  (denoted  $\hat{\gamma}$ ) and 95% confidence sets for this parameter, using respectively the asymptotic standard errors (2.27) (under the heading Wald-type), and the likelihood-ratio-type tests with i.i.d. Gaussian,  $t(\kappa)$  and normal mixture  $(\pi, \omega)$  errors, plus Gaussian GARCH errors (lower panel). <sup>10</sup> The values of  $\gamma$  in the Fieller-type confidence set are not rejected by the test defined in Theorem 3.1 to test  $\mathcal{H}(\gamma_0)$ . Rejection decisions are based on the largest MC p-values over all  $\kappa$  and  $(\pi, \omega)$ , respectively; we did not restrict maximization to the confidence set for these parameters here. As expected in view of the  $\mathcal{H}_B$  test results, the exact confidence sets are empty for several subperiods. The usefulness of the asymptotic confidence

<sup>10.</sup> Note that some values of  $\hat{\gamma}$  are high. Nonetheless, comparing the average real market return for those subperiods with our estimate of  $\gamma$  reveal that these high occurrences of  $\gamma$  are consistent with subperiods during which the estimated zero-beta rate is higher than the market portfolio return. This is an illustration of finding, ex post, a linear relationship between risk and return with a negative slope. Furthermore, rerunning our analysis using 10-year subperiods leads to  $\gamma$  estimates below the benchmark average return.

TABLE 6
QML-based point and set estimates for the zero-beta portfolio rate

Sample	$ar{R}_M$	$ar{r}_f$	$\hat{\gamma}$	Wald-type
$\overline{1927 - 30}$	0.0045	0.0045	0.0047	[-0.0037, 0.0130]
1931 - 35	0.0103	0.0025	-0.0130	[-0.0301, 0.0039]
1926 - 40	0.0031	-0.0006	-0.0069	[-0.0192, 0.0055]
1941 - 45	0.0097	-0.0042	0.0117	[0.0037, 0.0198]
1946 - 50	0.0021	-0.0051	-0.0219	[-0.0189, -0.0070]
1951 - 55	0.0145	0.0001	0.0024	[-0.0015, 0.0064]
1956 - 60	0.0086	0.0002	0.0156	[0.0109, 0.0202]
1961 - 65	0.0080	0.0014	0.0571	[0.0398, 0.0744]
1966 - 70	0.0008	0.0004	0.0169	[0.0096, 0.0242]
1971 - 75	-0.0061	-0.0010	0.0150	[0.0030, 0.0270]
1976 - 80	0.0056	-0.0012	-0.0096	[-0.0169, -0.0024]
1981 - 85	0.0081	0.0037	0.0197	[0.0125, 0.0268]
1986 - 90	0.0088	0.0020	0.0053	[-0.0024, 0.0131]
1991 - 95	0.0104	0.0011	0.0010	[-0.0130, 0.0062]

95% Confidence set, Fieller-type

Sample	Normal errors	Student t errors	Mixture errors	GARCH
1927 – 30	[-0.0133, 0.0227]	[-0.0143, 0.0229]	[-0.0141, 0.0227]	[-0.0125, .020]
1931 - 35	[-0.0509, 0.0225]	[-0.0520, 0.0225]	[-0.0157, 0.0227]	[-0.0517, .0217]
1926 - 40	[-0.0341, 0.0187]	[-0.0350, 0.0190]	[-0.0349, 0.0817]	[-0.0300, .0175]
1941 - 45	[-0.0045, 0.0275]	[-0.0048, 0.0287]	[-0.0045, 0.0283]	[-0.0025, .0275]
1946 - 50	Ø	Ø	Ø	Ø
1951 - 55	Ø	Ø	Ø	Ø
1956 - 60	Ø	[0.0149, 0.0161]	Ø	Ø
1961 - 65	Ø	Ø	Ø	Ø
1966 - 70	Ø	Ø	Ø	Ø
1971 - 75	[-0.0069, 0.0454]	[-0.0081, 0.0488]	[-0.0069, 0.0531]	[-0.0050, .0450]
1976 - 80	Ø	Ø	Ø	Ø
1981 - 85	[0.0059, 0.0371]	[0.0051, 0.0376]	[0.0051, 0.0387]	[0.0075, 0.0350]
1986 - 90	Ø	Ø	Ø	Ø
1991 - 95	[-0.0285, 0.0147]	[-0.0303, 0.0154]	[-0.0325, 0.0147]	[-0.0275, 0.0125]

Notes:  $\bar{R}_M$  is the average real market portfolio return for each subperiod,  $\bar{r}_f$  is the real average risk-free rate for each subperiod;  $\hat{\gamma}$  is the QML estimate of  $\gamma$ ; the remaining columns report 95% confidence sets for this parameter, using, respectively, the asymptotic standard errors  $(2.27)[\hat{\gamma}\pm 1.96\times \text{AsySE}(\hat{\gamma})]$ , the inverted test based on  $LR(\gamma_0)$  from Theorem 3.1, and the MC Gaussian p-value, the MMC p-value imposing multivariate  $t(\kappa)$  errors and mixture-of-normals  $(\pi,\omega)$  errors, and the MMC GARCH p-value. See Section 4 for details on the construction of these confidence sets. Non-Gaussian p-values are the largest MC p-values over the shape parameters  $\kappa$  or  $(\pi,\omega)$ . The GARCH p-value is the largest MC p-value over  $\phi_{1i},\phi_{2i}$  [from (2.17)]. Returns for the months of January and October 1987 are excluded from the data set

intervals is obviously questionable here. Other results which deserve notice are the empty sets for 1956–1960 subperiod; these sets correspond to the case where the efficiency bound test is significant (at 5%). <sup>11</sup>

To illustrate the differences between the asymptotic confidence set and ours, we next check whether the average real risk-free rate is contained in the confidence sets. For many subperiods, like 1966–1970, the evidence produced by the asymptotic and MC Fieller-type confidence intervals is similar. There are nonetheless cases where the set estimates do not lead to the same decision. For instance, for 1941–1945 and 1971–1975, the average risk-free rate is not included in the asymptotic confidence interval, while it is covered by our MMC confidence sets. These are

<sup>11.</sup> This can be checked by referring to Table 4: although the reported maximal p-values in this table are performed over the confidence set for  $\kappa$  and  $(\pi, \omega)$ , we have checked that the global maximal p-value leads to the same decision here.

	(1)	(2)	(3)	(4)
sample	$\tilde{\gamma} = \underset{\gamma_0}{\operatorname{argmin}} \mathcal{J}(\gamma_0)$	$\min_{\gamma_0} \mathscr{J}(\gamma_0)$	BND	95% Confidence set, MMC
1927 - 30	0.0090	71.29	0.650	[-0.0195, 0.0235]
1931 - 35	-0.0045	71.06	0.541	[-0.0240, 0.0250]
1926 - 40	-0.0045	54.52	0.620	[-0.0355, 0.0550]
1941 - 45	0.0415	163.26	0.143	[-0.0455, 0.0670]
1946 - 50	0.0000	133.76	0.121	[-0.0105, 0.0075]
1951 - 55	0.0075	104.93	0.250	[0.0000, 0.0120]
1956 - 60	0.0195	110.18	0.280	[-0.0385, 0.0415]
1961 - 65	0.0370	149.61	0.142	$[-0.0295, -0.0150] \cup [0.0250, 0.0670]$
1966 - 70	0.0090	168.54	0.081	[0.0045, 0.0135]
1971 - 75	0.0060	61.06	0.623	[-0.0180, 0.0067]
1976 - 80	0.0060	172.09	0.061	[-0.0225, 0.0135]
1981 - 85	0.0195	121.41	0.201	[0.0105, 0.0385]
1986 - 90	0.0030	184.38	0.030	Ø
1991 - 95	0.0100	53.60	0.841	$\{\le 0.0075\} \cup \{\ge 0.0310\}$

TABLE 7  $\mathcal{J}(\gamma_0)$  based inference on the zero-beta portfolio rate

Notes:  $\mathcal{J}(\gamma_0)$  is the HAC statistic in (2.28).  $\tilde{\gamma}$  is the minimum distance estimator from (2.29). Column (3) provides a bound MC p-value simulated at  $\tilde{\gamma}$  and maximized over  $\phi_{1i}, \phi_{2i}$  [from (2.17)]. Column (4) provides the confidence set for  $\gamma$  which inverts the inverted test based on  $\mathcal{J}(\gamma_0)$  and the MMC GARCH p-value; again, this is the largest MC p-value over  $\phi_{1i}, \phi_{2i}$  [from (2.17)]. Returns for the months of January and October 1987 are excluded from the data set. Given a 5% level, the cut-off of the BND p-value is 0.05; p-values that lead to significant tests with this benchmark are in bold. Note that the confidence set that inverts  $\mathcal{J}(\gamma_0)$  based on the asymptotic  $\chi^2(12)$  cut-off is empty for all subperiods.

cases where, using the asymptotic confidence interval, the hypothesis  $\gamma = r_f$  is rejected, whereas exact confidence sets indicate it should not be rejected. Conversely, in 1986–1990, the asymptotic confidence interval includes the average risk-free rate, whereas our confidence sets are empty.

In Table 7, we report the  $\mathcal{J}(\gamma_0)$  counterparts of the above QML-based tests (Columns 2 and 3) as well as point and set estimates of  $\gamma$  (Columns 1 and 4). Column (2) reports the values of our proposed J-test-type minimum  $\mathcal{J}(\gamma_0)$  statistic. In Column (3), MMC refers to the maximal MC p-value [over all  $(\phi_{1i}, \phi_{2i})$ ] for this statistic, assuming the GARCH specification (2.17), and the level is 5%; alternatively, an asymptotic  $\chi^2(12)$  critical value (21.03 for a 5% level) can be used. In column (1), we report the GMM-type point estimate (denoted  $\tilde{\gamma}$ ); the associated set estimate which inverts the  $\mathcal{J}(\gamma_0)$  MC Gaussian GARCH based test is reported in Column (4).

We first note that, on using the asymptotic critical value, a  $\mathcal{J}(\gamma_0)$  test would reject the model in all subperiods at level 5%. In contrast, the GARCH-MMC p-value is less than 5% only in the 1986–1990 subperiod. In view of our simulation results from Section 7, these results illustrate the serious implications of asymptotic test size distortions. Recall that the likelihood-ratio-based MC and MMC (Gaussian and non-Gaussian, with and without GARCH) tests reject the model at the 5% level in at least three other subperiods: 1946–1950, 1950–1955, 1960–1965. This reflects the test relative power, as illustrated in Section 7. Turning to the estimates of  $\gamma$ , we note that the  $\mathcal{J}(\gamma_0)$  based MMC confidence sets are substantially wider than the likelihood-ratio-based counterparts, only one confidence set is empty (in the 1986–1990 subperiods, in which case the model would be rejected), and the set is unbounded in the 1990–1995 subperiod. Had we relied on the asymptotic  $\chi^2(12)$  cut-off to invert the  $\mathcal{J}(\gamma_0)$  test, all confidence sets would be empty. Again, these observations line up with our simulation results.

The above procedures applied to the full data yields empty confidence sets using the exact GARCH corrected likelihood-ratio and  $\mathcal{J}(\gamma_0)$  criteria; the confidence interval using (2.27) is [0.0007,0.0088]. Since our subperiod analysis suggests that  $\gamma$  is temporally unstable, one must be careful in interpreting such results. On using a Bonferroni argument (that accounts for

time-varying parameters) based on the minimum (over subperiods) GARCH-corrected p-value that is 0.003 < 0.05/12, the model can be safely rejected at level 5%, over the full sample.

#### 9. CONCLUSION

This article proposes exact mean-variance efficiency tests when the zero-beta (or risk-free) rate is not observable, which raises identification difficulties. Proposed methods are robust to this problem as well as to portfolio repacking, and allow for heavy-tailed return distributions. In particular, a useful invariance result for MacKinlay and Richardson, 1991 HAC statistic is shown analytically. We also derive exact confidence sets for the zero-beta rate  $\gamma$ . While available Wald-type intervals are unreliable and lead to substantially different inference concerning  $\gamma$ , our confidence sets are valid in finite samples without assuming identification, and are empty by construction if efficiency is rejected.

We report a simulation study that illustrates the properties of our proposed procedures. Our results allow to disentangle small-sample problems from asymptotic failures arising from weak identification. We also examine efficiency of the market portfolio for monthly returns on NYSE CRSP portfolios. We find that efficiency is less rejected with non-normal assumptions. Exact confidence sets for  $\gamma$  differ importantly from asymptotic ones, and likelihood-ratio-based confidence sets are tighter than their Wald counterparts. All confidence sets nevertheless suggest that  $\gamma$  is not stable over time.

These results provide the motivation to extend our method to more general factor models, as discussed by (Campbell *et al.*, 1997, Chapter 6) and Shanken and Zhou (2007). These models raise the same statistical issues as the BCAPM, except that their definitional parameter is non-scalar. In this case, Fieller-type methods are clearly more challenging and raise worthy theoretical and empirical research questions.

#### A. APPENDIX: PROOFS

Proof of Theorem 3.1 Under (2.13) and  $\mathcal{H}(\gamma_0)$ , we have:  $T\hat{\Sigma} = \hat{U}'\hat{U} = K'W'M(X)WK$ ,  $T\hat{\Sigma}(\gamma_0) = K'W'\bar{M}(\gamma_0)WK$ . Then, under  $\mathcal{H}(\gamma_0)$ ,

$$\Lambda(\gamma_0) = \frac{|\hat{\Sigma}(\gamma_0)|}{|\hat{\Sigma}|} = \frac{|K^{'}W^{'}\bar{M}(\gamma_0)WK|}{|K^{'}W^{'}M(X)WK|} = \frac{|K^{'}|\left|W^{'}\bar{M}(\gamma_0)W\right||K|}{|K^{'}|\left|W^{'}M(X)W\right||K|} = \frac{\left|W^{'}\bar{M}(\gamma_0)W\right|}{|W^{'}M(X)W|},$$

hence  $P[LR(\gamma_0) \ge x] = P[T \ln(|W'\overline{M}(\gamma_0)W|/|W'M(X)W|) \ge x], \forall x.$ 

Proof of Theorem 3.2 For this model,

$$\hat{U} = M(X)U = M(X)WK = \hat{W}K, \quad \hat{W} = M(X)W.$$

If we denote by  $\hat{U}_t'$  and  $\hat{W}_t'$  the *t*-th rows of  $\hat{U}$  and  $\hat{W}$ , respectively, this means that  $\hat{U}_t = K'\hat{W}_t$ , for all *t*. Then,

$$\begin{split} \left(X_{t} \otimes \hat{U}_{t}\right) \left(X_{t-j} \otimes \hat{U}_{t-j}\right)' &= \left(X_{t} \otimes K' \hat{W}_{t}\right) \left(X_{t-j} \otimes K' \hat{W}_{t-j}\right)' = \left(X_{t} \otimes K' \hat{W}_{t}\right) \left(X'_{t-j} \otimes \hat{W}'_{t-j} K\right) \\ &= X_{t} X'_{t-j} \otimes K' \hat{W}_{t} \hat{W}'_{t-j} K = \left(I_{k} \otimes K'\right) \left(X_{t} \otimes \hat{W}_{t}\right) \left(X_{t-j} \otimes \hat{W}_{t-j}\right)' (I_{k} \otimes K) \end{split}$$

for  $j \ge 0, \dots$  Setting

$$\Psi_{j,T} = \frac{1}{T} \sum_{t=j+1}^{T} \left( X_{t} \otimes \hat{U}_{t} \right) \left( X_{t-j} \otimes \hat{U}_{t-j} \right)', S_{T} = \Psi_{0,T} + \sum_{j=1}^{q} \left( \frac{q-j}{q} \right) \left[ \Psi_{j,T} + \Psi'_{j,T} \right],$$

we see that

$$\Psi_{j,T} = \left(I_k \otimes K'\right) \hat{\Psi}_{j,T} \left(I_k \otimes K\right), \quad \hat{\Psi}_{j,T} = \frac{1}{T} \sum_{t=j+1}^T \left(X_t \otimes \hat{W}_t\right) \left(X_{t-j} \otimes \hat{W}_{t-j}\right)',$$

$$\begin{split} \Psi_{j,T} + \Psi_{j,T}^{'} &= \left(I_{k} \otimes K^{'}\right) \left[\hat{\Psi}_{j,T} + \hat{\Psi}_{j,T}^{'}\right] \left(I_{k} \otimes K\right), \\ S_{T} &= \left(I_{k} \otimes K^{'}\right) \hat{\Psi}_{0,T} \left(I_{k} \otimes K\right) + \sum_{j=1}^{q} \left(\frac{q-j}{q}\right) \left(I_{k} \otimes K^{'}\right) \left[\hat{\Psi}_{j,T} + \hat{\Psi}_{j,T}^{'}\right] \left(I_{k} \otimes K\right) = \left(I_{k} \otimes K^{'}\right) \hat{S}_{T} \left(I_{k} \otimes K\right), \\ \hat{S}_{T} &= \hat{\Psi}_{0,T} + \sum_{j=1}^{q} \left(\frac{q-j}{q}\right) \left[\hat{\Psi}_{j,T} + \hat{\Psi}_{j,T}^{'}\right]. \end{split}$$

On replacing  $S_T$  with  $(I_k \otimes K') \hat{S}_T (I_k \otimes K)$ , we get:

$$\left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] S_T \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] = \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K' \right] \hat{S}_T \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K \right].$$

Furthermore, since  $R(\gamma_0) = (1, \gamma_0) \otimes I_n = H(\gamma_0) \otimes I_n$  and  $H(\gamma_0)'$  is a k-dimensional vector in  $\tilde{\mathcal{H}}(\gamma_0) : R(\gamma_0)\vartheta = C'H(\gamma_0)' = 0$ , we see that:

$$R(\gamma_0)\hat{\vartheta} = \hat{C}'H(\gamma_0)',$$

$$R(\gamma_0) \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K' \right] = (H(\gamma_0) \otimes I_n) \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K' \right] = H(\gamma_0) \left( \frac{X'X}{T} \right)^{-1} \otimes K',$$

so  $\mathcal{J}(\gamma_0)$  can be rewritten as

$$\mathcal{J}(\gamma_0) = T \hat{\vartheta}' R(\gamma_0)' \left\{ R(\gamma_0) \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] S_T \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes I_n \right] R(\gamma_0)' \right\}^{-1} R(\gamma_0) \hat{\vartheta}$$

$$= T \hat{\vartheta}' R(\gamma_0)' \left\{ R(\gamma_0) \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K' \right] \hat{S}_T \left[ \left( \frac{X'X}{T} \right)^{-1} \otimes K \right] R(\gamma_0)' \right\}^{-1} R(\gamma_0) \hat{\vartheta}$$

$$= T H(\gamma_0) \hat{C} \mathcal{Q}(\gamma_0)^{-1} \hat{C}' H(\gamma_0)' = T H(\gamma_0) \hat{C} K^{-1} K \mathcal{Q}(\gamma_0)^{-1} K' \left( K' \right)^{-1} \hat{C}' H(\gamma_0)'$$

$$= T H(\gamma_0) \hat{C} K^{-1} \bar{\mathcal{Q}}(\gamma_0, W)^{-1} \left( K' \right)^{-1} \hat{C}' H(\gamma_0)'$$

where

$$\mathcal{Q}(\gamma_0) = \left\lceil H(\gamma_0) \left( \frac{X'X}{T} \right)^{-1} \otimes K' \right\rceil \hat{\mathsf{S}}_T \left\lceil \left( \frac{X'X}{T} \right)^{-1} H(\gamma_0)' \otimes K \right\rceil, \, \bar{\mathcal{Q}}(\gamma_0, W) = \left( K' \right)^{-1} \mathcal{Q}(\gamma_0) K^{-1}.$$

Under the null hypothesis  $H(\gamma_0)C = 0$ , we have:

$$H(\gamma_0)\hat{C} = H(\gamma_0)(X'X)^{-1}X'[XC+U] = H(\gamma_0)C + H(\gamma_0)(X'X)^{-1}X'U = H(\gamma_0)(X'X)^{-1}X'WK$$

$$\begin{split} \bar{\mathcal{Q}}(\gamma_0, W) &= \left(K'\right)^{-1} \left[ H(\gamma_0) \left(\frac{X'X}{T}\right)^{-1} \otimes K' \right] \hat{S}_T \left[ \left(\frac{X'X}{T}\right)^{-1} H(\gamma_0)' \otimes K \right] K^{-1} \\ &= \left[ H(\gamma_0) \left(\frac{X'X}{T}\right)^{-1} \otimes \left(K'\right)^{-1} K' \right] \hat{S}_T \left[ \left(\frac{X'X}{T}\right)^{-1} H(\gamma_0)' \otimes KK^{-1} \right] \\ &= \left[ H(\gamma_0) \left(\frac{X'X}{T}\right)^{-1} \otimes I_n \right] \hat{S}_T \left[ \left(\frac{X'X}{T}\right)^{-1} H(\gamma_0)' \otimes I_n \right] \end{split}$$

where we take into account the fact that  $H(\gamma_0)'$  is a column vector, hence

$$\mathcal{J}(\gamma_0) = TH(\gamma_0)(X'X)^{-1}X'WKK^{-1}\bar{\mathcal{Q}}(\gamma_0, W)^{-1}(K')^{-1}K'W'X(X'X)^{-1}H(\gamma_0)'$$

$$= TH(\gamma_0)(X'X)^{-1}X'W\bar{\mathcal{Q}}(\gamma_0, W)^{-1}W'X(X'X)^{-1}H(\gamma_0)'. \quad \|$$

Proof of Lemma 5.1 The Gaussian log-likelihood function for model (2.5) is

$$\ln[\tilde{L}(\tilde{Y},C,\Sigma)] = -\frac{T}{2}[n(2\pi) + \ln(|\Sigma|)] - \frac{1}{2}\operatorname{tr}[\Sigma^{-1}(\tilde{Y}-XC)'(\tilde{Y}-XC)] = \ln[L(Y,B,\Sigma)].$$

Setting  $\tilde{\Sigma}(C) \equiv \frac{1}{T}(\tilde{Y} - XC)'(\tilde{Y} - XC)$ , for any given value of C,  $\ln[\tilde{L}(\tilde{Y}, C, \Sigma)]$  is maximized by taking  $\Sigma = \tilde{\Sigma}(C)$  yielding the concentrated log-likelihood

$$\ln[\tilde{L}(\tilde{Y}, C, \Sigma)_c = -\frac{nT}{2}[(2\pi) + 1] - \frac{T}{2}\ln(|\tilde{\Sigma}(C)|). \tag{A.1}$$

The Gaussian MLE of C thus minimizes  $|\tilde{\Sigma}(C)|$  with respect to C. Let us denote by  $\hat{C}(Y)$  the unrestricted MLE of C so obtained, and by  $\hat{C}(Y;\gamma_0)$  and  $\hat{C}_B(Y)$  the restricted estimators subject to  $\hat{\mathscr{H}}(\gamma_0)$  and  $\hat{\mathscr{H}}_B$ , respectively. Suppose that  $\tilde{Y}$  is replaced by  $\tilde{Y}_* = \tilde{Y}A$  where A is a non-singular  $n \times n$  matrix. We need to show that  $LR_*(\gamma_0) = LR(\gamma_0)$  and  $LR_{B*} = LR_B$ , where  $LR_*(\gamma_0)$  and  $LR_{B*}$  represent the corresponding test statistics based on the transformed data. Following this transformation,  $|\tilde{\Sigma}(C)|$  becomes:

$$\begin{split} |\tilde{\Sigma}_{*}(C_{*})| &= \left| \frac{1}{T} (\tilde{Y}_{*} - XC_{*})' (\tilde{Y}_{*} - XC_{*}) \right| = \left| \frac{1}{T} A' (\tilde{Y} - XC_{*}A^{-1})' (\tilde{Y} - XC_{*}A^{-1})A \right| \\ &= |A'A| \left| \frac{1}{T} (\tilde{Y} - XC)' (\tilde{Y} - XC) \right| = |A'A| |\tilde{\Sigma}(C)| \end{split} \tag{A.2}$$

where  $C = C_*A^{-1}$ . Then  $|\tilde{\Sigma}(C_*)|$  is minimized by  $\hat{C}_*(Y_*) = \hat{C}(Y)A$  and  $|\tilde{\Sigma}_*(\hat{C}_*(Y_*))| = |A'A| |\tilde{\Sigma}(\hat{C}(Y))$ . On observing that  $H(\gamma_0)C = 0 \Longleftrightarrow H(\gamma_0)CA = 0 \Longleftrightarrow H(\gamma_0)C_* = 0$  for any  $\gamma_0$ , the restricted estimators of C under  $\tilde{\mathscr{H}}(\gamma_0)$  and  $\tilde{\mathscr{H}}_B$  are transformed in the same way:  $\hat{C}_*(Y_*; \gamma_0) = \hat{C}(Y; \gamma_0)A$  and  $\hat{C}_{*B}(Y_*) = \hat{C}_B(Y)A$ . This entails that  $|\tilde{\Sigma}_*(\hat{C}_*(Y_*; \gamma_0))| = |A'A| |\tilde{\Sigma}(\hat{C}(Y; \gamma_0))|$  and  $|\tilde{\Sigma}_*(\hat{C}_*(Y_*))| = |A'A| |\tilde{\Sigma}(\hat{C}_B(Y))|$ , so that

$$\tilde{\Lambda}_{*}(\gamma_{0}) = \frac{|\tilde{\Sigma}_{*}(\hat{C}_{*}(Y_{*};\gamma_{0}))|}{|\tilde{\Sigma}_{*}(\hat{C}_{*}(Y_{*}))|} = \frac{|\tilde{\Sigma}(\hat{C}(Y;\gamma_{0}))|}{|\tilde{\Sigma}(\hat{C}(Y))|} = \tilde{\Lambda}(\gamma_{0}), \tag{A.3}$$

$$\tilde{\Lambda}_{B*} = \frac{|\tilde{\Sigma}_*(\hat{C}_{*B}(Y_*))|}{|\tilde{\Sigma}_*(\hat{C}_*(Y_*))|} = \frac{|\tilde{\Sigma}(\hat{C}_B(Y))|}{|\tilde{\Sigma}(\hat{C}(Y))|} = \tilde{\Lambda}_B. \tag{A.4}$$

Finally, in view of (2.20) and (2.26), we have  $LR_*(\gamma_0) = T \ln[\tilde{\Lambda}_*(\gamma_0)] = T \ln[\tilde{\Lambda}(\gamma_0)] = LR(\gamma_0)$  and  $LR_{B*} = T \ln(\tilde{\Lambda}_{B*}) = T \ln(\tilde{\Lambda}) = LR_B$ .

*Proof of Theorem 5.2* Consider a transformation of the form  $\tilde{Y}_* = \tilde{Y}K^{-1}$  or, equivalently,  $Y_* = YK^{-1} + \tilde{R}_M \iota'_n (I - K^{-1})$ . Using (2.1) and (2.13), we then have:

$$Y_{*} = (XB + WK)K^{-1} + \tilde{R}_{M}\iota'_{n}(I - K^{-1}) = XBK^{-1} + \tilde{R}_{M}\iota'_{n}(I - K^{-1}) + W$$

$$= (\iota_{T}a' + \tilde{R}_{M}\beta')K^{-1} + \tilde{R}_{M}\iota'_{n}(I - K^{-1}) + W$$

$$= \tilde{R}_{M}\iota'_{n} + [\iota_{T}a' + \tilde{R}_{M}(\beta - \iota_{n})']K^{-1} + W$$

$$= \tilde{R}_{M}\iota'_{n} + X(B - \Delta)K^{-1} + W = \tilde{R}_{M}\iota'_{n} + X\bar{B} + W$$
(A.5)

where  $\bar{B} = (B - \Delta)K^{-1}$  and  $\Delta = [0, \iota_n]'$ . Using Lemma 5.1,  $LR(\gamma_0)$  and  $LR_B$  can be viewed as functions of  $Y_*$ , and depend on (B, K) only through  $\bar{B} = (B - \Delta)K^{-1}$ . Under  $\mathcal{H}_B$ , the nuisance parameter only involves  $\gamma$  and  $(\beta - \iota_n)'K^{-1}$ . Now the distribution of  $LR(\gamma_0)$  and  $LR_B$  can be explicitly characterized by using (A.3)–(A.4) and observing that

$$\begin{split} \tilde{\Lambda}(\gamma_0) &= \frac{|\tilde{\Sigma}_*\big(\hat{C}_*(Y_*;\gamma_0)\big)|}{|\tilde{\Sigma}_*\big(\hat{C}_*(Y_*)\big)|} = \frac{|\hat{W}(\gamma_0)'\hat{W}(\gamma_0)|}{|\hat{W}'\hat{W}|}, \\ \tilde{\Lambda}_B &= \frac{|\tilde{\Sigma}_*\big(\hat{C}_{*B}(Y_*)\big)|}{|\tilde{\Sigma}_*\big(\hat{C}_*(Y_*)\big)|} = \frac{\inf\left\{|\tilde{\Sigma}_*\big(\hat{C}_*(Y_*;\gamma_0)\big)|:\gamma_0 \in \Gamma\right\}}{|\tilde{\Sigma}_*\big(\hat{C}_*(Y_*)\big)|} = \inf\left\{\tilde{\Lambda}(\gamma_0):\gamma_0 \in \Gamma\right\}, \end{split}$$

where  $\hat{W}(\gamma_0) = \bar{M}(\gamma_0)(Y_* - \tilde{R}_M \iota_n') = \bar{M}(\gamma_0)(X\bar{B} + W) = \bar{M}(\gamma_0)\{[\iota_T a' + \tilde{R}_M (\beta - \iota_n)']K^{-1} + W\} = \bar{M}(\gamma_0)\{[\iota_T (a' + \gamma_0 (\beta - \iota_n)') + (\tilde{R}_M - \gamma_0 \iota_T)(\beta - \iota_n)']K^{-1} + W\} = \bar{M}(\gamma_0)\{\iota_T (a + \gamma_0 (\beta - \iota_n)')K^{-1} + W\} \text{ and } \hat{W} = M(X)W.$  Under  $\mathscr{H}_B$  where  $a = -\gamma(\beta - \iota_n)$ ,  $\hat{W}(\gamma_0) = (\gamma_0 - \gamma)\bar{M}(\gamma_0)\iota_T (\beta - \iota_n)'K^{-1} + \bar{M}(\gamma_0)W$ . The theorem then follows on observing that  $LR(\gamma_0) = T \ln[\tilde{\Lambda}(\gamma_0)]$  and  $LR_B = T \ln(\tilde{\Lambda}_B)$ . Further information can be drawn from the singular value decomposition of  $\bar{B}$ . Let r be the rank of  $\bar{B}$ . Since  $\bar{B}$  is a  $2 \times n$  matrix, we have  $0 \le r \le 2$  and we can write:

$$\bar{B} = PDQ', \quad D = [\bar{D}, 0], \quad \bar{D} = \text{diag}(\lambda_1^{1/2}, \lambda_2^{1/2}),$$
 (A.6)

where D is a  $2 \times n$  matrix,  $\lambda_1$  and  $\lambda_2$  are the two largest eigenvalues of  $\bar{B}'\bar{B}$  (where  $\lambda_1 \ge \lambda_2 \ge 0$ ),  $Q = [Q_1, Q_2]$  is an orthogonal  $n \times n$  matrix whose columns are eigenvectors of  $\bar{B}'\bar{B}$ ,  $Q_1$  is a  $2 \times r$  matrix that contains eigenvectors associated with the non-zero eigenvalues of  $\bar{B}'\bar{B}$ ,  $P = [P_1, P_2]$  is a  $2 \times 2$  orthogonal matrix such that  $P_1 = \bar{B}Q_1D_1^{-1}$  and  $D_1$  is a diagonal

matrix that contains the non-zero eigenvalues of  $\bar{B}'\bar{B}$ , setting  $P = P_1$  and  $D_1 = \bar{D}$  if r = 2, and  $P = P_2$  if r = 0; see (Harville, 1997, Theorem 21.12.1). Using Lemma 5.1 and Theorem 5.2,  $LR(\gamma_0)$  and  $LR_B$  may then be reexpressed as:

$$LR(\gamma_0) = T \ln \left( |\tilde{W}(\gamma_0)'\tilde{W}(\gamma_0)| / |\tilde{W}'\tilde{W}| \right), \quad LR_B = \inf \left\{ LR(\gamma_0) : \gamma_0 \in \Gamma \right\}, \tag{A.7}$$

$$\tilde{W} = \hat{W}Q = M(X)\bar{W}, \quad \bar{W} = WQ, \quad \tilde{W}(\gamma_0) = \hat{W}(\gamma_0)Q = \bar{M}(\gamma_0)(XPD + \bar{W}),$$
 (A.8)

 $PD = [P\bar{D}, 0]$  and  $P\bar{D}$  has at most 3 free coefficients (P is orthogonal). Under  $\mathcal{H}_B$ ,

$$\begin{split} \tilde{W}(\gamma_0) &= \bar{M}(\gamma_0) \iota_T \Big[ (\gamma_0 - \gamma) \big( \varphi' \varphi \big)^{1/2} \bar{\varphi}' \Big] + \bar{M}(\gamma_0) \bar{W}, \\ \varphi &= Q' \Big( K^{-1} \Big)' (\beta - \iota_n), \quad \bar{\varphi} &= \varphi / \big( \varphi' \varphi \big)^{1/2}. \end{split}$$

Define  $\Phi = [\bar{\varphi}, \bar{\Phi}]$  as an orthogonal matrix such that  $\Phi' \Phi = \Phi \Phi' = I_n$ , so

$$\Phi' \Phi = \begin{bmatrix} \bar{\varphi}' \bar{\varphi} & \bar{\varphi}' \bar{\Phi} \\ \bar{\Phi}' \bar{\varphi} & \bar{\Phi}' \bar{\Phi} \end{bmatrix} = \begin{bmatrix} 1 & 0 \\ 0 & I_{n-1} \end{bmatrix}, \qquad \bar{\varphi}' \Phi = \begin{bmatrix} 1 & 0 & \cdots & 0 \end{bmatrix}. \tag{A.9}$$

Then as in (A.7),  $LR(\gamma_0)$  and  $LR_B$  may again be expressed under  $\mathcal{H}_B$  as:

$$LR(\gamma_0) = T \ln \left( |\tilde{W}_B(\gamma_0)'\tilde{W}_B(\gamma_0)| / |\tilde{W}_B'\tilde{W}_B| \right), \quad LR_B = \inf \left\{ LR(\gamma_0) : \gamma_0 \in \Gamma \right\}, \tag{A.10}$$

$$\tilde{W}_B = \tilde{W}\bar{\Phi} = M(X)\bar{W}_B, \quad \bar{W}_B = \bar{W}\Phi, \tag{A.11}$$

$$\tilde{W}_B(\gamma_0) = \tilde{W}(\gamma_0)\bar{\Phi} = \bar{M}(\gamma_0)\iota_T \varphi_B' + \bar{M}(\gamma_0)\bar{W}_B, \tag{A.12}$$

where  $\varphi_B' = (\gamma_0 - \gamma) (\varphi' \varphi)^{1/2} \bar{\varphi}' \Phi = (\gamma_0 - \gamma) (\varphi' \varphi)^{1/2} [1 \ 0 \ \cdots \ 0]$  which involves at most one free coefficient. When W is non-Gaussian, the distributions of  $LR(\gamma_0)$  and  $LR_B$  may be influenced by  $\bar{B}$  through Q in  $\bar{W}$ . Under the Gaussian assumption (2.14), the rows of  $\bar{W}$  are i.i.d.  $N(0, I_n)$ , so that  $LR(\gamma_0)$  and  $LR_B$  follow distributions that depend on (B, K) only through  $P\bar{D}$ . Under  $\mathscr{H}_B$ , since the rows of  $\bar{W}_B$  are i.i.d.  $N(0, I_n)$ , this distribution involves only one nuisance parameter, in accordance with the result from (Zhou, 1991, Theorem 1), derived through a different method.

*Proof of Theorem 6.1*  $\mathcal{H}_B = \bigcup_{\gamma_0} \mathcal{H}(\gamma_0)$ . Since  $LR_B = \inf \{LR(\gamma_0) : \gamma_0 \in \Gamma\}$ , we have  $LR_B \leq LR(\gamma_0)$ , for any  $\gamma_0$ , hence  $P[LR_B \geq x] \leq P_{(B,K)}[LR(\gamma_0) \geq x]$ ,  $\forall x$ , for each  $\gamma_0$  and for any (B,K) compatible with  $\mathcal{H}(\gamma_0)$ . Furthermore, under  $\mathcal{H}_B$ , there is a value of  $\gamma_0$  such that the distribution of  $LR(\gamma_0)$  is given by Theorem 3.1, which entails (6.1). The result for the Gaussian special case then follows upon using (3.2).

*Proof of Theorem* 6.2 The result follows from (6.7), (3.12), and the inequalities  $\hat{p}_N^{\rm U}(LR_{\rm B};\gamma,\nu) \leq \hat{p}_N^{\rm U}(LR_{\rm B};\Gamma,\nu)$  and  $\hat{p}_N^{\rm U}(LR_{\rm B};\gamma,\nu) \leq \hat{p}_N^{\rm U}(LR_{\rm B};\gamma,\Omega_{\varnothing}) \leq \hat{p}_N^{\rm U}(LR_{\rm B};\Gamma,\Omega_{\varnothing})$ .

 $\begin{aligned} & \textit{Proof of Theorem 6.3} \quad \text{When $\nu$ is specified, by (6.6), (2.25), and (3.11), we have: } \hat{p}_N^U(LR_{\mathrm{B}}; \hat{\gamma}, \nu) \equiv p_N[LR_{\mathrm{B}}^{(0)}] \overline{LR}_N(\hat{\gamma}, \nu)] = \\ & p_N[LR^{(0)}(\hat{\gamma})] \overline{LR}_N(\hat{\gamma}, \nu)] = \hat{p}_N(LR; \hat{\gamma}, \nu), \text{ hence } \sup\{\hat{p}_N(LR; \gamma_0, \nu): \gamma_0 \in \Gamma\} \leq \alpha \Rightarrow \hat{p}_N(LR; \hat{\gamma}, \nu) \leq \alpha \Rightarrow \hat{p}_N^U(LR_{\mathrm{B}}; \hat{\gamma}, \nu) \leq \alpha \Rightarrow \sup\{\hat{p}_N(LR; \gamma_0, \nu): \gamma_0 \in \Gamma\} \leq \alpha \text{ means that } C_\gamma^{LR}(\alpha, \nu) \text{ is empty, } \hat{p}_N^U(LR_{\mathrm{B}}; \hat{\gamma}, \nu) > \alpha \Rightarrow \sup\{\hat{p}_N(LR; \gamma_0, \nu): \gamma_0 \in \Gamma\} > \alpha \Rightarrow C_\gamma^{LR}(\alpha, \nu) \neq \emptyset. \text{ For $\nu$ unknown,} \end{aligned}$ 

$$\begin{split} \hat{p}_{N}^{\mathrm{U}}(LR_{\mathrm{B}}; \hat{\gamma}, \Omega_{\mathcal{D}}) &= \sup\{\hat{p}_{N}^{\mathrm{U}}(LR_{\mathrm{B}}; \hat{\gamma}, \nu_{0}) : \nu_{0} \in \Omega_{\mathcal{D}}\} = \sup\{p_{N}[LR_{\mathrm{B}}^{(0)} \big| \overline{LR}_{N}(\hat{\gamma}, \nu_{0})] : \nu_{0} \in \Omega_{\mathcal{D}}\} \\ &= \sup\{p_{N}[LR^{(0)}(\hat{\gamma}) \big| \overline{LR}_{N}(\hat{\gamma}, \nu_{0})] : \nu_{0} \in \Omega_{\mathcal{D}}\} = \sup\{\hat{p}_{N}(LR; \hat{\gamma}, \nu) : \nu_{0} \in \Omega_{\mathcal{D}}\}, \end{split}$$

hence  $\sup\{\hat{p}_N(LR;\gamma_0,\nu_0):\gamma_0\in\Gamma,\nu_0\in\Omega_\varnothing\}\leq\alpha\Rightarrow\sup\{\hat{p}_N(LR;\hat{\gamma},\nu_0):\nu_0\in\Omega_\varnothing\}\leq\alpha\Rightarrow\hat{p}_N^U(LR_B;\hat{\gamma},\Omega_\varnothing)\leq\alpha$  and  $\hat{p}_N^U(LR_B;\hat{\gamma},\Omega_\varnothing)>\alpha\Rightarrow\sup\{\hat{p}_N(LR;\gamma_0,\nu_0):\gamma_0\in\Gamma,\nu_0\in\Omega_\varnothing\}>\alpha\Rightarrow\mathcal{C}_v^{LR}(\alpha;\varnothing)\neq\emptyset$ .

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