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TESTING FOR PARAMETER INSTABILITY AND STRUCTURAL CHANGE IN PERSISTENT PREDICTIVE REGRESSIONSˆ

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Testing for Parameter Instability and Structural Change in Persistent Predictive Regressionsˆ Torben G. Andersen and Rasmus T. Varneskov NBER Working Paper No. 28570 March 2021 JEL No. G12,G17

ABSTRACT

This paper develops parameter instability and structural change tests within predictive regressions for economic systems governed by persistent vector autoregressive dynamics. Specifically, in a setting where all – or a subset – of the variables may be fractionally integrated and the predictive relation may feature cointegration, we provide sup-Wald break tests that are constructed using the Local speCtruM (LCM) approach. The new tests cover both parameter variation and multiple structural changes with unknown break dates, and the number of breaks being known or unknown. We establish asymptotic limit theory for the tests, showing that it coincides with standard testing procedures. As a consequence, existing critical values for tied-down Bessel processes may be applied, without modification. We implement the new structural change tests to explore the stability of the fractionally cointegrating relation between implied- and realized volatility (IV and RV). Moreover, we assess the relative efficiency of IV forecasts against a challenging time-series benchmark constructed from high-frequency data. Unlike existing studies, we find evidence that the IV-RV cointegrating relation is unstable, and that carefully constructed time-series forecasts are more efficient than IV in capturing low-frequency movements in RV.

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1 Introduction

Predictive regressions are routinely invoked to assess the forecast prowess of candidate predictors for the future value of an economic quantity of interest. This approach provides great flexibility in accommodating multiple predictors, and it affords direct and transparent interpretability of the findings. Under standard regularity conditions, robust inference techniques may be applied to test for the significance of an individual regressor or the joint significance of (a subset of) the regressors.

The use of predictive regressions and least squares inference in time series contexts is, however, subject to controversy due to the inherent nature of economic and financial data, which has spurred a large literature dealing with potential pitfalls. One major issue is the strong persistence displayed by many economic time series. In particular, the spurious regression case, where one $I(1)$ process is projected onto another independent $I(1)$ process, generates large size distortions in standard least squares significance tests and artificially inflates model-fit measures such as the R^2 statistic; see e.g., Granger & Newbold (1974) and Phillips (1986). Similar problems arise for nearly integrated systems, e.g., Ng & Perron (1997) and Valkanov (2003), or if the variables are fractionally integrated, $I(d)$, where d is possibly a non-integer, see Tsay & Chung (2000) . Moreover, the regressions may be subject to sizable predictive biases if the system features endogenous correlation between the regression errors and innovations to the predictor variables, e.g., Stambaugh (1999) and Phillips & Lee (2013).

A second major complication is potential instability of the predictive regression relation. The evidence for the structural breaks in the data generating process is extensive within economics and finance; see, e.g., Stock & Watson (1996), Paye & Timmermann (2006), Peseran, Pettenuzzo & Timmermann (2006), Chen & Hong (2012), Dangl & Hailing (2012) and Farmer, Schmidt & Timmermann (2019). This issue is particularly salient when estimating predictive relations using data spanning multiple decades, and sometimes even centuries, or if using cointegration models to generate long-horizon forecasts, relying critically on accurate identification of the long-run equilibrium, see, e.g., Christoffersen & Diebold (1998). Since predictive regressions are reduced-form specifications, it is therefore not surprising that they evolve along with the economic environment over long time spans.

Several studies develop inference techniques for systems with persistent variables. For example, fractional (and standard) cointegration frameworks facilitate inference on linear predictive relations, whose error is purged of (at least some of) the persistence in the original variables; see, e.g., Robinson & Marinucci (2003), Christensen & Nielsen (2006), and Johansen & Nielsen (2012) for parametric and semi-parametric procedures. Moreover, robust techniques have been proposed for the nearly integrated case by, among others, Cavanagh, Elliott & Stock (1995), Jansson & Moreira (2006), and Elliott, Müller & Watson (2015). These procedures are reviewed by Phillips & Lee (2013), who develop the IVX methodology, allowing for stationary, nearly integrated, unit root and locally explosive processes. Finally, Andersen & Varneskov (2020a) provides a robust procedure for systems where all variables may be fractionally integrated of potentially different orders, covering both (asymptotically) stationary and non-stationary processes, and where the predictive relation may feature fractional cointegration, with a priori unknown persistence and cointegration properties. Furthermore, their Local speCtruM (LCM) approach accommodates regressor endogeneity and achieves asymptotic Gaussian inference.

Issues related to parameter instability and structural change have similarly received considerable attention. Andrews (1993), Andrews & Ploberger (1994) and Diebold & Chen (1996) develop tests for a structural change assuming that the break point is unknown, and Bai & Perron (1998) provide a comprehensive treatment for multiple structural changes with the timing, and possibly also the number, of breaks allowed to be unknown. These studies, however, require the system to be weakly dependent. In contrast, Hansen (1992), Hansen (2003) and Kerjiwal & Perron (2010), among others, develop similar structural change tests in $I(1)-I(0)$ cointegration models that also allow for $I(0)$ regressors. Moreover, Hidalgo & Robinson (1996) extend the sup-Wald test by Andrews (1993) to systems with stationary fractional integration, and Georgiev, Harvey, Leybourne & Taylor (2018) consider nearly integrated regressors. No prior study, however, has provided tests in the flexible persistence setting of Andersen & Varneskov $(2020a)$, that is, when the system can feature a mixture between stationary and non-stationary fractionally integrated variables and (fractional) cointegration, which is not known prior to the analysis. In fact, as discussed in the reviews by Perron (2006) and Casini & Perron (2019), existing procedures cannot simultaneously accommodate $I(0)$ and $I(1)$ errors, because the latter induce spurious inference. In this paper, we tackle these problems.

Specifically, we provide new tests for parameter instability and structural change, applicable for economic systems in which the variables (or a subset thereof) may be fractionally integrated of potentially different orders, and the predictive relation may feature cointegration, thereby nesting standard weakly dependent as well as cointegration settings. Our tests connect the LCM approach with the sup-Wald testing framework by Andrews (1993) and Bai & Perron (1998). Hence, we provide tests for structural changes with unknown break dates, where the number of breaks may be either known or unknown ex ante. These tests are generally applicable as parameter instability tests, despite being designed for the detection of structural breaks. We establish an asymptotic limit theory for the new tests. Importantly, by leveraging the asymptotic Gaussian distribution theory obtained via the LCM approach, we show that our tests, despite allowing for diverse and highly persistent systems, attain identical asymptotic distributions as the corresponding tests in Andrews (1993) and Bai & Perron (1998), who consider weakly dependent settings without cointegration. Hence, their critical values for tied-down Bessel processes may readily be applied in our context. Interestingly, due to the robustness properties of the LCM procedure, we also accommodate regressor endogeneity in the testing framework. Consequently, the new tests may be used as "off-line" diagnostics for a wide array of predictive systems, when one wants to retrospectively examine whether breaks occur in a given sample.¹

We illustrate the finite sample properties of our LCM-based sup-Wald testing procedure in settings that accommodate flexible, fractionally integrated persistence in the predictive relation, both in the absence and presence of cointegration. Despite the semi-parametric nature of our approach, the test

¹In contrast, if one wants to construct forecasts in real time, two alternatives include using rolling windows to implement the predictive regressions, e.g., Pesaran & Timmermann (2007) and Clark & McCracken (2009), or the explicit modeling of time-variation in the parameters, e.g., Peseran et al. (2006) and Dangl & Hailing (2012). None of their results, however, pertain to our flexible setting with fractionally integrated variables and, possibly, cointegration.

has excellent finite sample size and power properties for sample sizes typically encountered in financial and macroeconomic applications. Moreover, it is robust towards tuning parameter selections.

We use our new sup-Wald testing procedure to study the stability of the predictive relation between option implied and realized volatility (IV and RV), which has received considerable attention in empirical finance and financial econometrics. Specifically, using predictive regressions, this literature has been concerned with examining the bias and efficiency of IV as a predictor of future RV. After correcting for various methodological issues, consensus is converging towards the conclusion that IV forecasts are biased, but efficient predictors of future volatility; see, e.g., Christensen & Prabhala (1998), Jiang & Tian (2005), Chernov (2007), Chraoenwong, Jenwittayaroje & Low (2009), Taylor, Yadav & Zhang (2010) and Cheng & Fung (2012). These studies, however, all rely on inference procedure for weakly dependent processes, despite comprehensive evidence that return RV exhibits long memory; e.g., Andersen & Bollerslev (1997), Andersen, Bollerslev, Diebold & Ebens (2001), Andersen, Bollerslev, Diebold & Labys (2001, 2003), Corsi (2009) and Varneskov & Perron (2018).

The evidence of long-memory in IV and RV has inspired a number of studies to assess whether their predictive relation is fractionally cointegrated and whether IV is "long-run unbiased" for the persistent variation in RV. In particular, Bandi & Perron (2006), Christensen & Nielsen (2006), Nielsen (2007), Kellard, Dunis & Sarantis (2010) and Nielsen & Frederiksen (2011) provide solid support for these hypotheses. In addition, Bollerslev, Sizova & Tauchen (2011) and Bollerslev, Osterrieder, Sizova & Tauchen (2013), Osterrieder, Ventosa-Santaularia & Vera-Valdes (2019), Li, Izzeldin & Yao (2020) confirm the findings of fractional cointegration and proceed to exploit the equilibrium relation between IV and RV to estimate (components of) the volatility risk premium and forecast asset returns.

We revisit the long-run relation between IV and RV by first examining the relative efficiency of the long-run information in IV for RV. We compare it to forecasts from a long-memory volatility model exploiting high-frequency data, thereby providing a more challenging benchmark than historical volatility, which typically is used in the literature. Second, we examine the stability of the fractional cointegrating relations using our new LCM-based sup-Wald tests. Unlike existing studies, we find the IV-RV cointegrating relation to be unstable, and that carefully constructed time-series forecasts are more efficient than IV in capturing low-frequency movements in RV. These results are interesting from a volatility forecasting perspective, indicating a change in the low-frequency characteristics of RV over time. In addition, they have implications for asset pricing and return prediction, because our evidence is at odds with the presumed stability of the cointegration models, which are often used to estimate the volatility risk premium component associated with a long-run equilibrium and to forecast returns. In fact, we provide reduced-form evidence that the significant forecasting power of the volatility risk premium for returns materializes during a period surrounding the global financial crises of 2008-2009, where the low-frequency relation between IV and RV breaks down, with the former being completely spanned by our long-memory time-series forecasts. In contrast, during periods where a "normal" longrun equilibrium between IV and RV prevails, the volatility risk premium has no significant predictive power for returns. These findings provide a challenge for existing asset pricing models that stipulate

a stationary relation between RV, IV, the volatility risk premium and future returns.

The paper proceeds as follows. Section 2 describes the modeling framework, the underlying assumptions, and the competing parameter (in)stability hypotheses. Section 3 introduces the LCM procedure and derives the requisite partial-sample limit theory. Section 4 provides the new LCM-based sup-Wald tests and establishes their asymptotic properties. The simulation study is contained in Section 5, and Section 6 explores the IV-RV relation empirically. Section 7 concludes. The proofs of the theoretical results are relegated to the Appendix together with additional simulation results.

Throughout the paper, "∼" and " \approx " signify, respectively, that the ratio of the left- and right-handside of the relevant equation tends to one and a non-zero constant in the limit, element-wise, while we use $\|\cdot\|$ to denote the Frobenius matrix norm, ∘ the Hadamard product, $i = \sqrt{-1}$, and I_k the kdimensional identity matrix. Moreover, the notation " \rightarrow ", " $\stackrel{\mathbb{D}}{\rightarrow}$ " and " $\stackrel{\mathbb{P}}{\rightarrow}$ " describe the limit, convergence in distribution and convergence in probability. Finally, we denote by "⇒" weak convergence under the uniform metric on the space $D[0, 1]$, as defined by Pollard (1984, Chapter V).

2 Framework and Competing Hypotheses

First, to set the stage, let $(1 - L)^d$ denote a generic fractional filter,

$$
(1-L)^d = \sum_{i=0}^{\infty} \frac{\Gamma(i-d)}{\Gamma(i+1)\Gamma(-d)} L^i,
$$
\n(1)

with $\Gamma(\cdot)$ being the gamma function. Next, we assume observations are available for a $(k+1) \times 1$ vector of economic variables $z_t = (y_t, x'_{t-1})'$, which obeys a Type II fractional model,

$$
D(L)(z_t - \mu) = v_t 1_{\{t \ge 1\}}
$$
\n
$$
(2)
$$

where μ is a $(k+1) \times 1$ vector of nonrandom unknown finite numbers, either capturing the means or initial values of the variables, $v_t = (e_t, u'_{t-1})'$ is a weakly dependent vector process, and,

$$
\mathbf{D}(L) = \text{diag}[(1-L)^{d_1}, \ldots, (1-L)^{d_{k+1}}].
$$

In this setting, following Andersen & Varneskov (2020a), we define and study the predictive relation between y_t and x_{t-1} through their weakly dependent components. Specifically, we assume,

$$
e_t = \mathcal{B}'_t u_{t-1} + \eta_t^{(b)}, \qquad t = 1, \dots, n,
$$
\n(3)

where $\eta_t^{(b)} = (1 - L)^b \eta_t$ for some constant $b \geq 0$, $\eta_t \sim I(0)$, and the k-dimensional parameter vector, **B**_t, decomposes as $\mathcal{B}_t \equiv (\beta'_t, \delta')'$ with $\beta_t \in \mathbb{R}^{k_{\beta}}, \delta \in \mathbb{R}^{k_{\delta}}$ and $k = k_{\beta} + k_{\delta}$.

The main difference between the relation (3) and the corresponding assumption in Andersen & Varneskov (2020*a*) concerns the potential time variation of β_t . They consider estimation and inference for predictability under the standard invariant parameter vector assumption,

$$
\mathcal{H}_0: \ \ \beta_t = \beta, \qquad \text{for all} \quad t = 1, \dots, n. \tag{4}
$$

Importantly, even under the restriction, $\beta_t = \beta$, the observation equation (2) and functional form (3) encompass most fractionally integrated systems within the long memory literature, and they translate into a well-defined, *balanced* relation between y_t and x_{t-1} , despite the variables exhibiting distinct degrees of fractional integration. This is readily seen by combining the two equations, obtaining,

$$
y_t = a + \mathcal{B}'\mathcal{Q}(L)\,x_{t-1} + v_t, \qquad t = 1, \ldots, n,\tag{5}
$$

where $\mathcal{Q}(L) = \mathcal{D}_x(L)(1-L)^{-d_1}, \mathcal{D}_x(L)$ is the $k \times k$ lower-right submatrix of the fractional filter $\mathcal{D}(L)$, $a = \mu_y - \mathcal{B}'\mathcal{Q}(L)\mu_x$ for $\mu = (\mu_y, \mu'_x)'$ and $v_t = (1 - L)^{-d_1} \eta_t^{(b)} = (1 - L)^{b - d_1} \eta_t$. The filtering matrix $\mathcal{Q}(L)$ adjusts the persistence of the predictors so that regression balance is achieved in this implied relation between the observed variables. If the system is balanced a priori, then $\mathcal{Q}(L) = I_k$, and the adjustment is trivial. Its presence, however, ensures that the implied dynamics of the dependent variable, $y_t \sim I(d_1)$, are equivalent, both in the presence and absence of predictive power,

$$
\mathcal{B} = \mathbf{0} \qquad \text{and} \qquad \mathcal{B} \neq \mathbf{0}, \tag{6}
$$

respectively, and apply for settings with and without fractional cointegration;

- (i) $\mathcal{B} = 0$ and $b = 0$: x_{t-1} contains no predicitive power for y_t .
- (ii) $\mathcal{B} \neq 0$ and $b = 0$: x_{t-1} partially spans the persistent component(s) of y_t .
- (iii) $\mathcal{B} \neq 0$ and $b > 0$: x_{t-1} cointegrate with y_t , spanning its persistent component(s).

Andersen & Varneskov (2020*a*) show that most predictive long-memory systems, with and without cointegration, are nested in scenarios (i)-(iii), and they propose the Local SpectruM (LCM) procedure (described below) to draw inference on \mathcal{B} , whose validity holds uniformly across these settings, without prior knowledge of the persistence of the system, or whether it features cointegration. Hence, the LCM procedure overcomes problems with spurious inference in persistent systems, as documented by, among others, Granger & Newbold (1974), Phillips (1986) and Tsay & Chung (2000).

Their analysis, however, does rest critically on the null hypothesis, H_0 . In contrast, in this paper, we are concerned with testing whether \mathcal{H}_0 holds, allowing for either parameter instability or structural change in (a subset of) \mathcal{B}_t under the alternative. In line with previous studies offering instability tests in regression contexts, see, e.g., Andrews (1993) and Bai & Perron (1998) for one and multiple breaks, respectively, our framework nests pure and partial structural change models, when $k_{\delta} = 0$ and $0 < k_{\delta} < k$. Our new structural change tests within this flexible long-memory setting substantively generalizes the dynamic specifications assumed in prior work, both under the null hypothesis, as conveyed by scenarios (i)-(iii), and under the alternative. The specific alternatives are detailed next.

2.1 Alternative Hypotheses

First, define the set $\Phi \in (\epsilon, 1 - \epsilon)$, for some $\epsilon > 0$, let $\{\varphi_1 < \cdots < \varphi_q\} \in \Phi$ denote the sample break fractions, and fix $\varphi_0 = 0$ and $\varphi_{q+1} = 1$. In addition, we label the corresponding observation subsets $\mathcal{T}_g = \{t = \lfloor \varphi_g n \rfloor + 1, \ldots, \lfloor \varphi_{g+1} n \rfloor\}$ for the $q + 1$ regimes $g = 0, \ldots, q$ and denote the number of observations in \mathcal{T}_g by $n_g = n(\mathcal{T}_g)$. We will entertain multiple alternative hypotheses concerning structural change of the form,

$$
\mathcal{H}_{A1}: \ \beta_t = \beta_g, \qquad \text{for} \quad t \in \mathcal{T}_g, \quad g = 0, \dots, q \,, \tag{7}
$$

where the break dates are unknown, while q may be known or unknown, a priori. This alternative allows the predictive power of x_{t-1} to be characterized by distinct regimes of the form (i)-(iii). For example, x_{t-1} may partially span the persistence of y_t in some subsets of the sample, but not in others, thereby switching between scenarios (i) and (ii). Moreover, we allow for predictive systems, where the cointegration properties for a subset of the regressors may break down in specific economic regimes, as long as the system remains in scenario (iii). To see this, suppose we have a cointegration model, where $y_t \in I(1),$ $\mathbf{x}_{t-1} = (x_{1,t-1}, x_{2,t-1})' \in I(1)$ and $b = 1$. Then, our setting may feature a partial structural change model, with $k_{\beta} = k_{\delta} = 1$, where cointegration arises through a time-invariant relation with $x_{2,t-1}$, but the predictive power of $x_{1,t-1}$ pertains only to certain regimes.

As a final example, the alternative hypothesis accommodates fractionally cointegrated systems, where persistence is balanced, that is, with $y_t \in I(d)$, $x_{t-1} \in I(d)$ and $0 \leq b \leq d$, and where the relative efficiency of (a subset of) the predictors may change over the course of the sample.

The flexibility of scenarios (i)-(iii), allowing all variables to have different fractional integration orders and the predictive relation to feature cointegration, in conjunction with the alternative hypothesis (7), facilitates testing for parameter changes in the vast majority of the predictive long-memory systems studied in the literature. Furthermore, it is robust to a feature not accommodated by prior structural break tests; namely, we remain agnostic as to whether the system involves cointegration. Specifically, Andrews (1993) and Bai & Perron (1998), among others, provide structural change tests with unknown break dates in weakly dependent settings, and Hansen (1992), Hansen (2003) and Kerjiwal & Perron (2010) consider cointegrated systems that may also feature $I(0)$ regressors, but their respective procedures cannot simultaneously accommodate $I(0)$ and $I(1)$ errors, because the latter induce spurious inference; see, e.g., the discussions in Perron (2006) and Casini & Perron (2019). In our setting, the errors may be $I(0)$ or $I(1)$, and we also accommodate fractional cointegration within the system. As will be explained in detail subsequently, this added generality stems from our use of the LCM procedure in the construction of the structural change tests.

Finally, we emphasize that, despite our tests being designed with alternatives such as the structural break specification (7) in mind, they will also have power against other alternatives of the form,

$$
\mathcal{H}_{A2}: \ \beta_t \neq \beta_s, \qquad \text{for some} \quad t,s \geq 1,
$$

as noted by Andrews (1993) in a weakly dependent setting without cointegration. Following the analysis in the latter, we show below that our new tests, indeed, have power against local alternatives of the form $\beta_t = \beta + \eta(t/n)/S_n(b)$ for some bounded function $\eta(t/n)$ on [0,1], that is not almost everywhere constant on Φ , and where $\mathcal{S}_n(b)$ is an increasing deterministic function of the sample size n, which depends on the cointegration strength, b. Hence, our tests will capture more general forms of parameter instability under the predictive scenarios (i)-(iii).

2.2 Assumptions

The theoretical analysis rests on formal assumptions for the components of equations (2) and (3). Moreover, we assume an estimator of the fractional integration orders, \hat{d}_i for $i = 1, ..., k+1$, satisfying mild consistency requirements, is available. The assumptions, stated in terms of $q_t = (u'_{t-1}, \eta_t)'$, mirror the corresponding ones in Andersen & Varneskov (2020a), and they are consistent with the regularity conditions underpinning the fractional cointegration analyses in, e.g., Robinson & Marinucci (2003), Christensen & Nielsen (2006) and Christensen & Varneskov (2017).

Assumption D1. The vector process q_t , $t = 1, \ldots$, is covariance stationary with spectral density matrix satisfying $f_{qq}(\lambda) \sim G_{qq}$ as $\lambda \to 0^+$, where the upper left $k \times k$ submatrix, G_{uu} , has full rank, and the $(k + 1)$ th element of the diagonal, $G_{\eta\eta}$, is strictly greater than zero. Moreover, there exists a $\omega \in (0,2]$ such that $|\mathbf{f}_{qq}(\lambda) - \mathbf{G}_{qq}| = O(\lambda^{\omega})$ as $\lambda \to 0^+$. Finally, let $\mathbf{G}_{qq}(i,k+1)$ be the $(i,k+1)$ th element of G_{qq} , which has $G_{qq}(i, k+1) = G_{qq}(k+1, i) = 0$ for all $i = 1, \ldots, k$.

Assumption D2. q_t is a linear process, $q_t = \sum_{j=0}^{\infty} A_j \epsilon_{t-j}$, with square summable coefficients $\sum_{j=0}^{\infty} ||A_j||^2 < \infty$, the innovations satisfy, almost surely, $\mathbb{E}[\epsilon_t|\mathcal{F}_{t-1}] = 0$ and $\mathbb{E}[\epsilon_t \epsilon'_t|\mathcal{F}_{t-1}] = I_{k+1}$, and the matrices $\mathbb{E}[\epsilon_t \otimes \epsilon_t \epsilon'_t | \mathcal{F}_{t-1}]$ and $\mathbb{E}[\epsilon_t \epsilon'_t \otimes \epsilon_t \epsilon'_t | \mathcal{F}_{t-1}]$ are nonstochastic, finite, and do not depend on t, with $\mathcal{F}_t = \sigma(\epsilon_s, s \leq t)$. There exists a random variable ζ such that $\mathbb{E}[\zeta^2] < \infty$ and, for all c and some C, $\mathbb{P}[\Vert q_t \Vert > c] \leq C \mathbb{P}[\vert \zeta \vert > c]$. For future reference, we denote the periodogram of ϵ_t by $J(\lambda)$.

Assumption D3. For $A(\lambda, i)$, the *i*-th row of $A(\lambda) = \sum_{j=0}^{\infty} A_j e^{ij\lambda}$, its partial derivative satisfies $\|\partial \mathbf{A}(\lambda, i)/\partial \lambda\| = O(\lambda^{-1} \|\mathbf{A}(\lambda, i)\|)$ as $\lambda \to 0^+$, for $i = 1, \ldots, k + 1$.

Assumption M. Let $0 \leq d_i < 2$ for all $i = 1, ..., k + 1$ and define $\underline{d} = \min_{i=1,...,k+1} d_i$. Moreover, suppose that the cointegration strength parameter satisfies $0 \leq b \leq \min(\underline{d}, 1)$.

Assumption F. Let $m_d \n\approx n^{\varrho}$ be a sequence of integers where $0 \leq \varrho \leq 1$, then, for all $i = 1, \ldots, k+1$ elements of z_t , we assume to have an estimator with the property,

$$
\widehat{d}_i - d_i = O_p\left(1/\sqrt{m_d}\right), \quad \text{and we then let,} \quad \widehat{D}(L) = \text{diag}\left[(1-L)^{\widehat{d}_1}, \ldots, (1-L)^{\widehat{d}_{k+1}}\right].
$$

The regularity conditions on q_t in Assumptions D1-D3 are standard in the literatures analyzing semi-parametric fractional cointegration as well as the estimation of multivariate fractional time series models, e.g., Shimotsu (2007) and Nielsen (2015). Specifically, conditions D1 and D3 impose a mild rate

of convergence for the spectral density $f_{qq}(\lambda)$ as $\lambda \to 0^+$, which depends on the smoothness parameter $\pi \in (0, 2]$. Moreover, D1 requires full rank of u_{t-1} and the vector being locally exogenous to η_t as $\lambda \to 0^+$, but not global exogeneity. That is, u_{t-1} and η_t may have non-trivial covariance for frequencies $\lambda \to c > 0$. Finally, D2 specifies linearity, martingale and moment conditions for q_t , allowing for general multivariate dependence, but rules out time-variation in the conditional covariance between the innovations as well as their third and fourth conditional moments. Assumption M restricts the fractional integration order of the variables to be below 2 as well as the cointegration range in relation to the memory of the system. The first part is innocuous, since it is satisfied by most macroeconomic and financial time series. Moreover, the cointegration restriction is rather intuitive. If the system is balanced $d_i = d, i = 1, \ldots, k + 1$, then we accommodate both $I(1)$ - $I(0)$ and fractional cointegration settings, without any additional restrictions. If the memory parameters differ, Assumption M implies that the cointegration relation must be balanced and, at least, weakly dependent. To see this, suppose $y_t \in I(0.3)$ and $x_{t-1} \in I(0.6)$. Then, the upper bound $b = 0.3$, in conjunction with the balanced regression requirement (5) , implies a cointegrating relation with $I(0)$ residuals. On the other hand, if $y_t \in I(0.6)$ and $x_{t-1} \in I(0.3)$, then the restriction generates $I(0.3)$ residuals. Hence, the regressors cannot purge y_t of more persistence than their own degree of fractional integration.

These conditions allow for z_t to constitute vector ARFIMA processes (for $\varpi = 2$), thus nesting VAR and integrated VAR dynamics as special cases with $d_i = 0$ and $d_i = 1$, respectively, for $i = 1, \ldots, k+1$, and $b = 0$. Generally, however, the properties of the variables depend on their integration orders. Specifically, the *i*th variable have (asymptotically) stationary long memory, when $0 < d_i < 1/2$, it is non-stationary when $d_i \geq 1/2$, and it has a well-defined mean for $d_i < 1$. Hence, the assumptions allow for very flexible persistence of the variables as well as fractional cointegration amongst them.

Finally, the conditions on d_i in Assumptions F are mild, essentially only requiring the existence of an estimator which, under the Assumptions D1-D3 and M, is consistent. Examples of estimators, that satisfy these requirements for the entire range, $0 \leq d_i < 2$, are the parametric ARFIMA estimators by Hualde & Robinson (2011) and Nielsen (2015), the mean-robust exact Local Whittle (ELW) estimator by Shimotsu (2010) and the trimmed ELW (TELW) estimator by Andersen & Varneskov (2020a).

Remark 1. The LCM procedure, and thus our new tests below, further accommodates regressor endogenenity in the spirit of Stambaugh (1999) and Pastor & Stambaugh (2009) as well as regressors that have been pre-estimated as fractional cointegration residuals. These effects are formally analyzed in Andersen & Varneskov (2020a, Section 4), who show that robustness is achieved by slightly stricter trimming conditions than those required by Assumption T below. As these conditions are not binding in our empirical analysis, we refrain from formally introducing endogeity and pre-estimated regressors here for simplicity of exposition, noting, however, that all our subsequent qualitative results will apply.

3 Partial Sample LCM Inference

This section introduces a partial sample version of the two-step LCM estimator by Andersen & Varneskov (2020a). Furthermore, it provides a joint functional limit theory for $q + 1$ partitions of the sample under the null hypothesis of parameter stability and develop associated feasible inference procedures. These estimators, and their resulting limit theory, are of independent interest but, importantly, also form the basic building blocks in the design of the structural change tests in Section 4.

3.1 Partial Sample LCM Estimation and Limit Theory

The LCM estimation procedure is carried out in two steps.

Step 1: Fractional filtering. As in Assumption F, we use a consistent estimator of the fractional integration orders d_i , for $i = 1, ..., k + 1$, to compute $D(L)$, and then fractionally filter z_t to obtain estimates of v_t , the weakly dependent components of the variables in equation (2) ,

$$
\widehat{\boldsymbol{v}}_t \equiv (\widehat{e}_t, \widehat{\boldsymbol{u}}_{t-1}')' = \widehat{\boldsymbol{D}}(L)\boldsymbol{z}_t, \tag{8}
$$

without, however, accounting for the mean, or initial value, of z_t . Rather than treating "de-meaning" of the series on a case-by-case basis, depending on the magnitude of d_i , we account for the residual impact of the mean component, $\hat{D}(L) \mu$, in a unified manner during the second stage estimation.

Step 2: Medium Band Least Squares. For a given partition $g = 0, \ldots, q$, denote the Fourier frequencies for observation set \mathcal{T}_g by $\lambda_j(g) = 2\pi j/n_g$, $j = 1, \ldots, n_g$. Then, the partial sample discrete Fourier transform and cross-periodogram are defined as,

$$
\boldsymbol{w}_h(\lambda_j,g) = \frac{1}{\sqrt{2\pi n_g}} \sum_{t \in \mathcal{T}_g} \boldsymbol{h}_t \, e^{\mathrm{i}t\lambda_j(g)}, \qquad \boldsymbol{I}_{hk}(\lambda_j,g) = \boldsymbol{w}_h(\lambda_j,g) \, \overline{\boldsymbol{w}}_k(\lambda_j,g), \tag{9}
$$

respectively, where h_t and k_t are generic (and compatible) vector time series. Moreover, let $I_{hk}(\lambda_j, g)$ = $\Re(I_{hk}(\lambda_j, g)) + i\Im(I_{hk}(\lambda_j, g))$ be the real and imaginary decomposition of $I_{hk}(\lambda_j, g)$. Finally, define the corresponding trimmed discretely averaged co-periodogram (TDAC) as,

$$
\widehat{F}_{hk}(\ell,m,g) = \frac{2\pi}{n_g} \sum_{j=\ell_g}^{m_g} \Re(\mathbf{I}_{hk}(\lambda_j,g)), \qquad 1 \le \ell_g \le m_g \le n_g,
$$
\n(10)

where $\ell_g = \ell(n_g)$ and $m_g = m(n_g)$ are partition-specific trimming and bandwidth functions. Then, we can write the TDAC of $\hat{\mathbf{u}}_{t-1}$ as $\mathbf{F}_{\hat{u}\hat{u}}(\ell, m, g)$ and, similarly, of $\hat{\mathbf{u}}_{t-1}$ and \hat{e}_t as $\mathbf{F}_{\hat{u}\hat{e}}(\ell, m, g)$, and use these to define the partial sample medium band least squares (MBLS) estimator as,

$$
\widehat{\mathcal{B}}(\ell,m,g) = \widehat{F}_{\widehat{u}\widehat{u}}(\ell,m,g)^{-1} \widehat{F}_{\widehat{u}\widehat{e}}(\ell,m,g), \tag{11}
$$

for which $\ell_g, m_g \to \infty$ and $\ell_g/m_g + m_g/n_g \to 0$, as $n_g \to \infty$.² Moreover, by definition of the partial sample observation scheme, we have $n_g/n \to \varphi_{g+1} - \varphi_g > 0$ as $n \to \infty$, that is, each subsample grows proportionally to n . Importantly, the MBLS estimator in equation (11) is equivalent to the second-step estimator in Andersen & Varneskov (2020a), although it applies only to a subsample rather than the full set of observations.³ Hence, it inherits some useful properties. First, it is semiparametric about the dynamics of z_t , requiring only some specific structure on the spectrum as $\lambda \to 0^+$. Second, the trimming and bandwidth functions aid in the recovery of the parameter vector within each subsample by asymptotically eliminating the first-stage estimation errors through filtering. Specifically, the trimming component eliminates any slippage from the means, or initial values, $\hat{\bm{D}}(L)\bm{\mu}$, occurring at lower frequencies and, in conjunction with the bandwidth, the estimation errors given in Assumptions F.

Consistent with subsample estimation of the parameter vector \mathcal{B} , we require additional conditions on the partition-specific bandwidth and trimming functions. These are partial sample analogs of the corresponding assumptions in Andersen & Varneskov (2020a).

Assumption T. Let the bandwidth m^g n κ ^g and `^g n ν ^g with 0 < ν < κ < % ≤ 1 for sample partitions $g = 0, \ldots, q$. Moreover, recall that the parameter $\varpi \in (0, 2]$ measures the smoothness of the spectral density in Assumption D1. Then, the following cross-restrictions are imposed,

$$
\frac{m_g^{1+2\varpi}}{n_g^{2\varpi}} + \frac{\ell_g^{1+\varpi+b}}{n_g^{\varpi} \, m_g^{1/2+b}} + \frac{n_g^{1/2+b}}{m_d^{1/2} \, m_g^b \, \ell_g} + \frac{n_g^{1-2\underline{d}+b}}{m_g^{1/2-2\underline{d}+b} \, \ell_g^2} + \frac{n_g^b}{m_g^{1/2+b}} \to 0, \qquad \text{as} \quad n \to \infty.
$$

The conditions in Assumption T are mild. The first condition is standard for semiparametric estimation in the frequency domain, e.g., Robinson (1995) and Lobato (1999), while the last is needed, because we only require local exogeneity between u_{t-1} and η_t as $\lambda \to 0^+$. For the empirically relevant vector ARFIMA process ($\varpi = 2$), this condition implies $\kappa < 4/5$. The remaining conditions are LCMspecific upper and lower bounds in the trimming and bandwidth rates, ν and κ , which are discussed at length in Andersen & Varneskov (2020a). They depend on \underline{d} , b and $\underline{\rho}$, thus reflecting the persistence properties of the economic system, cointegration and the convergence rate of the first-step estimator of the fractional integration orders, respectively. For the vector ARFIMA process, the most important implications of conditions 2-4 of Assumption T amount to,

$$
(3/5 - (2\underline{d} - b)/5)/2 \vee (1 - \varrho)/2 + b(1 - \kappa) < \nu < \kappa < 4/5,\tag{12}
$$

²We have suppressed dependence on the (lagged) time t indicator in $\hat{F}_{\hat{u}\hat{u}}(\ell, m, g)$ and $\hat{F}_{\hat{u}\hat{e}}(\ell, m, g)$ to ease exposition. The dependence on time, however, is explicated when necessary, e.g., when establishing Theorem 1 below.

³The MBLS estimator is also related to the estimator in Christensen & Varneskov (2017), who, however, input observable variables rather than fractionally filtered ones. This has important implications; the framework in the latter only applies to (asymptotically) stationary processes and cannot handle situations where the integration orders differ across variables in the system. Moreover, the asymptotic analysis in Christensen & Varneskov (2017) assumes cointegration. However, since the latter shows robustness of their estimator (via trimming) towards outliers, structural breaks (in means), Markov switching means, certain deterministic trends, etc., which are known to contaminate co-periodograms at frequency ordinates close to the origin, see, e.g., Diebold & Inoue (2001) and Perron & Qu (2010), this suggests that our estimator will enjoy similar robustness properties. We do not formally analyze those effects here.

for the trimming rate, thereby generating a conservative bound $3/10 < \nu$ for standard choices of the parameters ϱ and κ . The last condition generates a mild lower bound restriction on the bandwidth rate, $b/(1/2 + b) \le \min(d, 1)/(1/2 + \min(d, 1)) < \kappa$. Naturally, the trimming and bandwidth rates can be made subsample specific, as long as they satisfy Assumption T in each partition. We refrain from doing so for simplicity of exposition and to avoid dealing with a large set of tuning parameters.

Before stating the joint functional limit theory for the subsample LCM estimators, we introduce some auxiliary notation. Specifically, we let,

$$
\boldsymbol{\vartheta}(b) = \text{diag}\left[(\varphi_1 - \varphi_0)^{2b(\kappa - 1) - \kappa - 1}, \dots, (\varphi_{q+1} - \varphi_q)^{2b(\kappa - 1) - \kappa - 1} \right] \tag{13}
$$

be a $(q + 1)$ -dimensional diagonal matrix, depending on the sample partition, and,

$$
C(b) = G_{uu}^{-1} V(b) G_{uu}^{-1}, \quad \text{where} \quad V(b) = G_{uu} \frac{G_{\eta\eta}}{2(1+2b)}, \quad (14)
$$

denote the $k \times k$ asymptotic covariance matrix of the LCM estimator. Moreover, write the full sample Fourier frequencies as $\lambda_j = 2\pi j/n$, $j = 1, ..., n$, and let $\ell = \ell(n)$ and $m = m(n)$ be the full sample trimming and bandwidth sequences, which also satisfy Assumption T. Finally, let $Y(\varphi)$, for some $\varphi \in [0,1]$, denote a $k \times 1$ vector of independent Brownian motions on [0, 1], and we then define,

$$
\boldsymbol{Y}(\boldsymbol{\varphi}) = \begin{pmatrix} \boldsymbol{Y}(\varphi_1) - \boldsymbol{Y}(\varphi_0) \\ \vdots \\ \boldsymbol{Y}(\varphi_{q+1}) - \boldsymbol{Y}(\varphi_q) \end{pmatrix},
$$
\n(15)

with dimension $(q+1) \times 1$, for the sample partitions φ on the set Φ . We are now ready to characterize the joint limit theory for subsample LCM estimators given any partition, $g = 0, \ldots, q$.

Theorem 1. Let Assumptions D1-D3, F, M and T hold. Moreover, suppose that the mutual consistency condition $\max(0, (1 - 3\kappa/2)/(1 + \kappa/2)) < \infty \leq 2$ is satisfied, then, for every sequence of sample partitions $\boldsymbol{\varphi} = (\varphi_1, \ldots, \varphi_q)'$, whose elements are in the set Φ , it follows,

$$
\sqrt{m}\,\lambda_m^{-b}\,\begin{pmatrix} \widehat{\mathcal{B}}(\ell,m,0)-\mathcal{B}\\\vdots\\ \widehat{\mathcal{B}}(\ell,m,q)-\mathcal{B} \end{pmatrix}\quad\Longrightarrow\quad \left(\boldsymbol{\vartheta}(b)^{1/2}\circ\left(\boldsymbol{G}_{uu}^{-1}\boldsymbol{V}(b)^{1/2}\right)\right)\boldsymbol{Y}(\boldsymbol{\varphi})\,.
$$

Theorem 1 extends Andersen & Varneskov $(2020a,$ Theorem 1) to allow for non-overlapping partitions of the sample and demonstrates that subsample LCM estimators enjoy a joint Gaussian uniform limit. Importantly, the convergence rate depends on whether $\mathcal{B} \neq 0$ and, if this is the case, whether the system features cointegration. Specifically, the rate is \sqrt{m} in its absence $(b = 0)$, consistent with semiparametric estimators in the frequency domain, e.g., Brillinger (1981, Chapters 7-8). Conversely, the rate is $\sqrt{m}\lambda_m^{-b} \asymp \sqrt{m}(n/m)^b$ in its presence $(b > 0)$, in analogy with super consistency properties in

standard cointegration settings. Moreover, cointegration lowers the asymptotic variance, as reflected by the scale $1/(2(1+2b))$. The joint Gaussian uniform distribution theory is remarkable; holding in scenarios with and without cointegration, across (asymptotically) stationary and non-stationary variables in the observable persistent system (2), and for both weak $(b < 1/2)$ and strong $(b \ge 1/2)$ cointegration settings. This feature is exactly what allows our structural change tests to remain, a priori, agnostic about the persistence of the system and whether it features cointegration, thus considerably generalizing the scope relative to, e.g., Andrews (1993), Bai & Perron (1998) and Kerjiwal & Perron (2010), who either rely on standard \sqrt{n} asymptotic properties of OLS (or GMM) estimators in weakly dependent settings or rate- n results in cointegration settings.

3.2 Feasible Inference

The design of our structural change tests below require estimates of the asymptotic covariance matrix for each sample partition. To this end, we must provide consistent subsample estimators of the longrun covariance matrix G_{uu} and variance $G_{\eta\eta}$. The main challenge is that we observe \hat{v}_t , not v_t raising issues similar to those that motivated us to apply the robust MBLS estimator in the second step of the LCM procedure. The residuals η_t are latent, and we estimate them as,

$$
\widehat{\eta}_t^{(b)}(g) = \widehat{e}_t - \widehat{\mathcal{B}}(\ell, m, g)' \widehat{\mathbf{u}}_{t-1}, \qquad \widehat{\eta}_t(g) \equiv (1-L)^{-\widehat{b}(g)} \widehat{\eta}_t^{(b)}(g), \quad t \in \mathcal{T}_g,
$$
\n(16)

for each subsample $g = 0, \ldots, q$, where $\hat{b}(g)$ denotes a consistent estimator of b. The intuition behind the estimator $\hat{\eta}_t(g)$ is the following; we must undo the (over-)differencing, indicated by $b \neq 0$, to recover estimates of the residuals, that are then used to estimate $G_{\eta\eta}$. This is carried out on each subsample for our estimates to be valid under the null, \mathcal{H}_0 , and the alternative hypothesis, \mathcal{H}_{A1} .

Next, we use \hat{v}_t and $\hat{\eta}_t(g), g = 0, \ldots, q$, to define a generic class of subsample estimators of the asymptotic covariance matrices,

$$
\widehat{G}_{hh}(\ell_G, m_G, g) = \frac{1}{m_{G,g} - \ell_{G,g} + 1} \sum_{j=\ell_{G,g}}^{m_{G,g}} \Re(\mathbf{I}_{hh}(\lambda_j, g)), \qquad (17)
$$

for some arbitrary vector h_t , where $m_{G,g} = m_G(n_g)$ and $\ell_{G,g} = \ell_G(n_g)$ are other, partition-specific, bandwidth and trimming functions, which, as for MBLS, are used to eliminate estimation errors from fractional filtering, low-frequency slippage from $\hat{D}(L)\mu$, as well as estimation errors from both MBLS and $b(g)$. Now, using equation (17), the asymptotic covariance of $\mathcal{B}(\ell, m, g)$ is estimated as,

$$
\widehat{\text{AVAR}}(g, b) = \widehat{G}_{\widehat{u}\widehat{u}}(\ell_G, m_G, g)^{-1} \frac{\widehat{G}_{\widehat{\eta}\widehat{\eta}}(\ell_G, m_G, g) \,\lambda_{m_g}^{2\widehat{b}(g)}(g)}{2\left(1 + 2\widehat{b}(g)\right) m_g},\tag{18}
$$

which implicitly accounts for the scale $(\varphi_{g+1} - \varphi_g)^{2b(\kappa-1)-\kappa}$ by applying $\lambda_{m_g}(g)$ and m_g instead of the full sample versions λ_m and m. To retain flexibility in terms of implementing the inference and testing

procedures, in analogy to our treatment of the estimators d_i , $i = 1, ..., k + 1$, we assume an estimator $b(g)$ is available, for $g = 0, \ldots, q$, which satisfies mild consistency requirements. For example, these conditions are satisfied by the TELW estimator of Andersen & Varneskov $(2020a)$. Before proceeding, we also need to impose restrictions on the new bandwidth and trimming parameters.

Assumption B. Let $m_{b,g} \n\approx n_g^{\varepsilon}$, with $g = 0, \ldots, q$, be a partition-specific sequence of integers where, for some $0 < \varepsilon \le 1$, the subsample estimators satisfy $\hat{b}(g) - b = O_p(1/\sqrt{m_{b,g}})$.

Assumption T-G. Let $m_{G,g} \approx n_g^{\kappa_G}$ and $\ell_{G,g} \approx n_g^{\nu_G}$ with $0 < \nu_G < \kappa_G \le \varrho \le 1$, for $g = 0, \ldots, q$, and define the sequence of integers $m_g = m_d \wedge m_{b,g} \wedge m_g$, then the following cross-restrictions are imposed on the trimming and bandwidth parameters: $n/(m_{G,g} \ell_{G,g}^2) + n^2/(m_{G,g} \ell_{G,g}^2 \underline{m}_g) \to 0$, as $n \to \infty$.

Finally, to illustrate the use of the subsample inference, we consider general linear hypotheses on the parameters **B** as $\mathcal{RB} = r$ for some $h \times k$ selection matrix **R** and $h \times 1$ vector **r**. The following theorem establishes validity of the feasible inference and partition-specific testing.

Theorem 2. Let the conditions of Theorem 1 and Assumptions B and T-G hold. Then, for every sequence of partitions $\varphi = (\varphi_1, \ldots, \varphi_q)'$, whose elements are on the set Φ , it follows,

(a)
$$
||m_g \lambda_{m_g}^{-2b}(g) \widehat{\text{AVAR}}(g, b) - C(b)|| \xrightarrow{\mathbb{P}} 0
$$
 for all $g = 0, ..., q$, and,
\n(b) $\mathcal{LCM}(\ell, m, g) \equiv (\mathcal{R}\widehat{\mathcal{B}}(\ell, m, g) - r)' (\mathcal{R}\widehat{\text{AVAR}}(g, b) \mathcal{R}')^{-1} (\mathcal{R}\widehat{\mathcal{B}}(\ell, m, g) - r) \xrightarrow{\mathbb{D}} \chi_h^2$.

Theorem 2 shows that Wald test statistics may be constructed for every sample partition and, importantly, have standard χ^2 -distribution theory. This feature is essential for the design and validity of our structural break tests, which are introduced next.

4 Testing for Parameter Instability and Structural Change

First, we provide a new test for general parameter instability as well as a single structural break with unknown break date, exploiting the Gaussian asymptotic limit theory for the LCM approach along with the sup-Wald testing procedure of Andrews (1993). Second, we extend this test to better accommodate multiple structural changes, occurring at unknown time points, where the number of regimes may be known or unknown, a priori, leveraging insights from Bai & Perron (1998).

4.1 Testing for Parameter Instability and a Single Structural Change

We exploit the standard asymptotic properties of the subsample LCM estimators in Theorems 1 and 2 to construct our Wald test statistic for a *single* structural break, $g = 1$,

$$
\mathcal{W}_n(\varphi_1) \equiv \left(\widehat{\boldsymbol{\beta}}(\ell,m,0) - \widehat{\boldsymbol{\beta}}(\ell,m,1) \right)^{\prime} \left(\widehat{\mathbf{AVAR}}_{\beta}(0,b) + \widehat{\mathbf{AVAR}}_{\beta}(1,b) \right)^{-1} \left(\widehat{\boldsymbol{\beta}}(\ell,m,0) - \widehat{\boldsymbol{\beta}}(\ell,m,1) \right),
$$

in line with the alternative hypothesis \mathcal{H}_{A1} , where $\widehat{\beta}(\ell, m, g)$ denotes the first k_{β} elements of the subsample LCM estimate $\hat{\mathcal{B}}(\ell, m, g)$ and, similarly, $\widehat{\mathbf{AVAR}}_{\beta}(g, b)$ denotes the upper-left $k_{\beta} \times k_{\beta}$ submatrix of $\mathbf{AVAR}(q, b)$. Then, we introduce a new class of sup-Wald tests,

$$
\mathcal{P}_n(1) \equiv \sup_{\varphi_1 \in \Phi} \mathcal{W}_n(\varphi_1), \tag{19}
$$

with, again, $\Phi \in (\epsilon, 1 - \epsilon)$ for some $0 < \epsilon < 1$. The asymptotic properties of the $\mathcal{P}_n(1)$ class of parameter instability tests under the null hypothesis are formalized by the following theorem.

Theorem 3. Suppose the conditions of Theorems 1-2 hold. Let $\zeta = \kappa/2 + (1 - \kappa)b$, and,

$$
Q_{k_{\beta}}(\varphi_1,\zeta) = \frac{\widetilde{\mathbf{Y}}_{k_{\beta}}(\varphi_1,\zeta)^{\prime}\widetilde{\mathbf{Y}}_{k_{\beta}}(\varphi_1,\zeta)}{((1-\varphi_1)^{2\zeta}+\varphi_1^{2\zeta})(\varphi_1(1-\varphi_1))}, \quad \text{where}
$$

$$
\widetilde{\mathbf{Y}}_{k_{\beta}}(\varphi_1,\zeta) = \left(\mathbf{Y}_{k_{\beta}}(\varphi_1)(1-\varphi_1)^{\zeta+1/2} - (\mathbf{Y}_{k_{\beta}}(1)-\mathbf{Y}_{k_{\beta}}(\varphi_1))\varphi_1^{\zeta+1/2}\right),
$$

and with $Y_{k_{\beta}}(\varphi_1)$ containing the first k_{β} elements of the vector of independent Brownian motions, $\mathbf{Y}(\varphi_1)$. Then, the following convergence results hold for all $\varphi_1 \in \Phi$,

$$
\mathcal{W}_n(\varphi_1) \Rightarrow Q_{k_{\beta}}(\varphi_1,\zeta), \qquad \mathcal{P}_n(1) \stackrel{\mathbb{D}}{\rightarrow} \sup_{\varphi_1 \in \Phi} Q_{k_{\beta}}(\varphi_1,\zeta).
$$

Despite the challenging setting, where all variables may exhibit different degrees of fractional integration and the system may feature cointegration, as conveyed by scenarios in (i)-(iii), the LCM-based sup-Wald test, $\mathcal{P}_n(1)$, enjoys asymptotic properties under the null hypothesis, which are identical to those achieved by the corresponding test in Andrews (1993) under weak dependence. This readily follows as,

$$
Q_{k_{\beta}}(\varphi_1,\zeta) \triangleq \frac{\left(\boldsymbol{Y}_{k_{\beta}}(\varphi_1) - \boldsymbol{Y}_{k_{\beta}}(1)\varphi_1\right)' \left(\boldsymbol{Y}_{k_{\beta}}(\varphi_1) - \boldsymbol{Y}_{k_{\beta}}(1)\varphi_1\right)}{\varphi_1\left(1 - \varphi_1\right)}\,,\tag{20}
$$

which is the tied-down Bessel process describing the asymptotic distribution theory in Andrews (1993, Theo. 3). The standardization in the construction of $\mathcal{W}_n(\varphi_1)$, thus, eliminates any potential changes in the asymptotic distribution of the LCM-based sup-Wald test stemming from the, possibly, nonstandard convergence rate of LCM, as signified by ζ , where $\zeta = 1/2$ corresponds to \sqrt{n} . Consequently, our test is readily implementable in practice, with critical values provided in Andrews (1993, Table 1). Moreover, one may apply the identical arguments, as for Theorems 4-5 and Corollary 2 of the latter, due to the LCM Gaussian limit theory, to show that $\mathcal{P}_n(1)$ has local power against alternatives of the form \mathcal{H}_{A2} . To obtain equivalent results, we simply set $\mathcal{S}_n(b) = \sqrt{m}\lambda_n^{-b}$ instead of \sqrt{n} , in line with the differences in convergence rates between LCM and GMM in their respective settings.

Remark 2. Due to the standard asymptotic behavior of the Wald test statistics, $W_n(\varphi_1)$, one may also consider exp-Wald and mean-Wald tests in the spirit of Andrews $\mathcal{C}_{\mathcal{A}}$ Ploberger (1994), using the former as building blocks. However, Kim & Perron (2009) show that tests for structural change based on either sup-Wald or exp-Wald formulations compares favorably to mean-Wald tests as well as tests based on Langrange Multiplier (LM) and Likelihood Ratio (LR) measures, according to an approximate Bahadur measure of efficiency. Hence, we abstain from pursuing such extensions here.

4.2 Testing for Multiple Structual Changes with Unknown Break Points

The LCM-based sup-Wald test in equation (19) is asymptotically valid against alternative hypotheses involving multiple structural changes, as shown Andrews (1993, Theo. 4-5, Cor. 2). Its finite sample power may be lacking, however; see, e.g., Vogelsang (1997). To alleviate concerns about the latter, we amend our LCM approach with the tests for multiple structural changes in Bai & Perron (1998), designed in weakly dependent settings and absent cointegration, to explicitly account for $(q+1)$ regimes with unknown break dates. To this end, let us define,

$$
\mathcal{R}(\widehat{\beta}, q) \equiv \left(\widehat{\beta}(\ell, m, 0)' - \widehat{\beta}(\ell, m, 1)', \dots, \widehat{\beta}(\ell, m, q - 1)' - \widehat{\beta}(\ell, m, q)'\right)', \tag{21}
$$

$$
\mathcal{A}(\widehat{\beta},q) \equiv \text{diag}\left(\widehat{\text{AVAR}}_{\beta}(0,b) + \widehat{\text{AVAR}}_{\beta}(1,b), \ldots, \widehat{\text{AVAR}}_{\beta}(q-1,b) + \widehat{\text{AVAR}}_{\beta}(q,b)\right), \quad (22)
$$

of dimensions $q k_{\beta} \times 1$ and $q k_{\beta} \times q k_{\beta}$, respectively. Then, we apply our subsample LCM framework and generalize the Wald statistic and sup-Wald test in equation (19) as,

$$
\mathcal{P}_n(q) \equiv \sup_{\boldsymbol{\varphi} \in \Phi} \mathcal{W}_n(\boldsymbol{\varphi}), \qquad \mathcal{W}_n(\boldsymbol{\varphi}) = \boldsymbol{\mathcal{R}}(\boldsymbol{\hat{\beta}}, q) \boldsymbol{\mathcal{A}}(\boldsymbol{\hat{\beta}}, q)^{-1} \boldsymbol{\mathcal{R}}(\boldsymbol{\hat{\beta}}, q) / (q k_{\boldsymbol{\beta}}), \qquad (23)
$$

This class of tests requires specifying the number of structural changes under the alternative, which may not be desirable in some applications. Hence, as also considered by Bai & Perron (1998), we further introduce a "double maximum" version of our LCM-based sup-Wald test,

$$
\mathcal{D}_n(q_{\max}) \equiv \max_{1 \le q \le q_{\max}} \mathcal{P}_n(q), \tag{24}
$$

for some upper bound q_{max} . The main advantage of "double maximum" tests, as thoroughly discussed by Bai & Perron (2006) in a weakly dependent setting, is that they alleviate problems with nonmonotonic power, if the true number of breaks is larger than what was implemented in the testing procedure, e.g., if we implement $P_n(1)$ when $q = 2$. Moreover, the power of the former is almost as high as what can be achieved by $\mathcal{P}_n(q)$ in cases where q is known.

Next, as for our sup-Wald test in equation (19), the properties of the $\mathcal{P}_n(q)$ and $\mathcal{D}_n(q_{\text{max}})$ classes of structural change tests are characterized under the null hypothesis in the following theorem.

Theorem 4. Suppose the conditions of Theorems 1-2 hold. Let $\zeta = \kappa/2 + (1 - \kappa)b$ and

$$
Q_{k_{\beta}}(\varphi,\zeta,g) = \frac{\widetilde{\boldsymbol{Y}}_{k_{\beta}}(\varphi,\zeta,g)'\widetilde{\boldsymbol{Y}}_{k_{\beta}}(\varphi,\zeta,g)}{((\varphi_{g+1}-\varphi_{g})^{2\zeta}+(\varphi_{g+2}-\varphi_{g+1})^{2\zeta}})\left((\varphi_{g+1}-\varphi_{g})(\varphi_{g+2}-\varphi_{g+1})\right)}, \text{ where}
$$

$$
\widetilde{\boldsymbol{Y}}_{k_{\beta}}(\varphi,\zeta,g) = \left(\Delta \boldsymbol{Y}_{k_{\beta}}(\varphi_{g+1})(\varphi_{g+2}-\varphi_{g+1})^{\zeta+1/2}-\Delta \boldsymbol{Y}_{k_{\beta}}(\varphi_{g+2})(\varphi_{g+1}-\varphi_{g})^{\zeta+1/2}\right),
$$

 $for g=0,\ldots,q-1,$ and with $\Delta Y_{k_{\beta}}(\varphi_{g+1})=Y_{k_{\beta}}(\varphi_{g+1})-Y_{k_{\beta}}(\varphi_g),$ with $Y_{k_{\beta}}(\varphi_g)$ containing the first k_{β} elements of the vector of independent Brownian motions, $\mathbf{Y}(\varphi_g)$. Moreover, define,

$$
\mathcal{Q}_{k_{\beta}}(\boldsymbol{\varphi}, \zeta,q) \; = \; \frac{1}{q k_{\beta}} \, \sum_{g=0}^{q-1} Q_{k_{\beta}}(\boldsymbol{\varphi}, \zeta,g) \, .
$$

Then, the following convergence results hold, for all $\varphi \in \Phi$,

$$
\mathcal{P}_n(q) \overset{\mathbb{D}}{\to} \sup_{\boldsymbol{\varphi} \in \Phi} \mathcal{Q}_{k_\beta}(\boldsymbol{\varphi}, \zeta, q) \,, \qquad \mathcal{D}_n(q_{max}) \overset{\mathbb{D}}{\to} \max_{1 \leq q \leq q_{max}} \sup_{\boldsymbol{\varphi} \in \Phi} \mathcal{Q}_{k_\beta}(\boldsymbol{\varphi}, \zeta, q) \,.
$$

Theorem 4 shows that the LCM-based sup-Wald tests for multiple structural breaks with unknown break dates deliver an asymptotic distribution theory, which is equivalent to the corresponding limits in Bai & Perron (1998, Proposition 6), despite our more challenging setting. The latter is readily seen by each partition satisfying distributional equivalence with a multi-break extension of representation (20). Hence, we may apply the critical values in Bai & Perron (1998, Table 1) to implement our LCM-based tests. Moreover, the power of the latter against alternatives of the form \mathcal{H}_{A2} follow by the inequalities $\mathcal{P}_n(1) \leq 2\mathcal{P}_n(2) \leq q\mathcal{P}_n(q)$, and consistency of our test in equation (19).

The results in Theorems 3 and 4 significantly broaden the applicability of sup-Wald tests for structural changes and parameter instability. Contrary to prior studies, such as Andrews (1993), Hansen (1992), Bai & Perron (1998), Hansen (2003), and Kerjiwal & Perron (2010), we accommodate the following dynamic features: (i) all variables of the system may be fractionally integrated of potentially different orders; (ii) the predictive relation may be fractionally cointegrated, covering $I(1)$ - $I(0)$ cointegration as a special case; (iii) we remain agnostic about the persistence and cointegration properties, ex ante.⁴ These empirically relevant extensions are rendered feasible by our utilization of LCM, and its Gaussian limit theory, in the design of the tests.

Remark 3. An interesting alternative to applying the critical values in Andrews (1993) and Bai \mathcal{B} Perron (1998) for structural change testing is to use bootstrap techniques such as those in Diebold $\&$ Chen (1996) and Hansen (2000), who document favorable size properties of their tests. Moreover, Georgiev et al. (2018) and Boldea, Cornea-Madeira \mathfrak{G} Hall (2019) show that bootstrap procedures are critical for valid testing of parameter instability in OLS regressions, if the predictors are persistent, modeled as local-to-unity processes, and endogenous, respectively. In contrast, we accommodate persistence through reliance on the flexible class of fractionally integrated processes, cointegration and endogenous regressors (cf., Remark 1). The primary distinction of our approach from theirs is that they correct biases in the asymptotic theory for OLS in such settings, whereas we eliminate corresponding biases via the LCM procedure, rendering our asymptotic theory Gaussian.

⁴See also the recent review in Casini & Perron (2019) for a discussion of the state of the literature.

5 Simulation Study

This section complements the asymptotic analysis by exploring the finite sample properties of the LCM-based sup-Wald test for parameter instability, $\mathcal{P}_n(1)$, through realistic simulations. Specifically, Section 5.1 describes the simulation setup, which is designed to mimic quantitative features of our empirical application in Section 6. Whereas Section 5.2 and 5.3 provide the main numerical results, we have included additional tables in Appendix B to supplement the analysis

5.1 Simulation Setting

We consider testing for parameter instability in an setting reminiscent of the ones in Hong (1996), Shao (2009) and Andersen & Varneskov (2020a), albeit allowing for time-varying predictive power of the regressors. In particular, we restrict attention to the "pure" structural break model, $\mathcal{B}_t = \beta_t$, and consider the univariate regressor case, thus writing x_{t-1} and β_t for the predictor and regression coefficient, respectively. Then, we generate fractional $ARMA(1, 0)$ processes as,

$$
(1 - L)^{d_1}(y_t - \mu_y) = \beta_t u_{t-1} + (1 - L)^b \eta_t, \quad (1 - L)^{d_2}(x_{t-1} - \mu_x) = u_{t-1}, \tag{25}
$$

where η_t and u_{t-1} are AR(1) processes with persistence parameters $\phi_{\eta} = \phi_u = 0.2$ and mean-zero Gaussian errors with volatility $\sigma_{\eta} = 0.5$ and $\sigma_{u} = 1$, respectively. Moreover, we fix the means $\mu_y = \mu_x = 1/2$, the integration orders $d_1 = d_2 = d$ and consider selections $d \in \{0.30, 0.60, 1.00\}$ as well as models with $(b = d)$ and without $(b = 0)$ fractional cointegration. These configurations nest long memory regressions with variables that are (asymptotically) stationary, non-stationary and it covers standard unit root regressions and cointegration models. For each specification, we consider sample sizes $n \in \{350, 700\}$, noting that the case $d = b = 0.60$ and $n = 350$ corresponds well with our empirical application to the IV-RV forecast relation in Section 6.

To analyze the size and power properties of $\mathcal{P}_n(1)$, we simulate the system (25) under the null hypothesis without breaks in the predictive relation $\mathcal{H}_0 : \beta_t = \beta = 1$ as well as examine various alternatives. For specificity, the alternative hypotheses stipulate a single break in the regression coefficient, that is, we consider parameters of the form $\beta_t(\varphi_1) = \beta_0 \mathbf{1}_{\{t \in \mathcal{T}_0(\varphi_1)\}} + \beta_1 \mathbf{1}_{\{t \in \mathcal{T}_1(\varphi_1)\}}$, for different values of the break fraction, φ_1 , and regime-dependent coefficients, β_0 and β_1 . Our baseline set of alternatives feature $\beta_0 = 0.50, \beta_1 = 1.50$ as well as $\mathcal{H}_1 : \varphi_1 = 0.25, \mathcal{H}_2 : \varphi_1 = 0.50$, and $\mathcal{H}_3 : \varphi_1 = 0.75$, thus capturing sizable breaks in the beginning, middle and end of the sample. In addition, we consider an equivalent set of *small* break alternatives, where $\beta_0 = 0.90$, $\beta_1 = 1.10$ and \mathcal{H}_4 : $\varphi_1 = 0.25$, \mathcal{H}_5 : $\varphi_1 = 0.50$ and \mathcal{H}_6 : $\varphi_1 = 0.75$, thus providing a very challenging set of tests in finite samples.⁵

We need to select an estimator of the fractional integration order along with a few tuning parameters

⁵In addition to the single break models, we have considered three different alternatives. First, $\beta_t = 0.50 \times \mathbf{1}_{\{t \leq 0.25n\}} +$ $1_{\{[0.25n] < t \leq [0.75n]\}} + 1.50 \times 1_{\{t > [0.75n]\}}$ has two structural breaks. Second, $\beta_t = 0.50(1 - t/n) + 1.50(t/n)$ captures a trending regression coefficient. Finally, we have examined a predictive relation with regression parameter exhibiting periodic fluctuation of the form $\beta_t = 1 + \frac{\sin(t/(n/10))}{n}$. The power results are indistinguishable from those reported for the baseline set of alternatives and, thus, omitted for brevity.

to implement the LCM-based sup-Wald test. First, we carry out the fractional filtering and estimation of the cointegration strength parameter using the TELW estimator of Andersen & Varneskov (2020a), as it applies to all of our testing scenarios (i.e., values of d and b).⁶ To this end, we follow the latter and use trimming $\ell_d = \lfloor n^{0.20} \rfloor$ and a bandwidth $m_d = \lfloor n^{0.75} \rfloor$. In addition, we will initially use the tuning parameters $\nu = \nu_G = 0.20$, $\kappa = 0.70$ and $\kappa_G = 0.90$ for all configurations of the predictive model. To verify that these selections, indeed, correspond well with the restrictions implied by Assumption T and equation (12), we have explicated the requisite bounds on the trimming and bandwidth rates for the different values of d and b in Table 1. Moreover, we examine to what extend the tuning parameter selections are robust in Section 5.3. Finally, we set $\epsilon = 0.20$, controlling the length of the search space for the sup-Wald test, Φ , the nominal size of the test to 5% (again, see Andrews (1993, Table 1) for critical values) and consider simulations with 1000 replications.

5.2 Simulation Results

The size and power results for the baseline alternatives are reported in Table 2 and deserve a few comments. First, the size of the test in the absence of cointegration $(b = 0)$ is 6.9-8.5%, thus close to the nominal 5% level. Moreover, it improves for the cointegration case $(b = d)$ when $d \leq 0.60$ due to a faster rate of convergence of the LCM estimator (cf. Theorem 1), whereas the test is slightly conservative for unit root persistence, $d = 1.00$. Second, the test is very powerful, as evidenced by the 100% rejection rates for all baseline alternatives. The corresponding rejection rates for the small break alternatives in Table 3 underscores its considerable finite sample power, which improves when the sample size is increased from $n = 350$ to $n = 700$. These results further show that the test has higher power to detect breaks in the middle of the sample than in the beginning or the end, for which, however, the rejection rates remain high. Finally, the simulations show that the power is generally higher for cointegrating regressions, not surprisingly, as these induce a faster rate of convergence.

The supplementary tables in Appendix B show that the full sample LCM estimate converges to a weighted average of the regime-dependent coefficients under the break alternative and that breaks generates higher variance of the coefficient estimates. Moreover, they demonstrate that the LCMbased sup-Wald procedure precisely estimates the break fraction for the baseline alternatives, yet have a bias for the small break alternatives, consistent with the power properties of the test.

5.3 Tuning Parameter Robustness

To assess the robustness of the LCM-based sup-Wald test for parameter instability, $\mathcal{P}_n(1)$, with respect to its tuning parameters, we consider various combinations with $\nu \in \{0.10, 0.20, 0.30\}, \ \kappa \in$ $\{0.50, 0.60, 0.70, 0.75\}$ and $\kappa_G \in \{0.75, 0.80, 0.90\}$, while fixing the remaining TELW configurations. Moreover, in order keep the exposition simple, we focus on predictive regressions where $d = 0.60$ and $n = 350$, which, as mentioned above, is closely aligned with our empirical application below, and we

 6 Note that estimation of the integration order and fractional filtering are performed using the full sample of observations before implementing the sup-Wald testing procedure on different partitions of the observations.

consider settings with $(b = d)$ and without $(b = 0)$ cointegration. The results are provided in Tables 4 and 5 for the *baseline* and *small* break alternatives, respectively. Whereas the properties of the sup-Wald testing procedure is robust to changes in the trimming rate, ν , as well as the bandwidth of the long-run covariance estimator, κ_G , we find the test properties to display some sensitivity to the MBLS bandwidth parameter, κ . Specifically, when $\kappa = 0.50$, we observe both size distortions and lower power of the procedure. The finite sample properties improve for $\kappa = 0.60$, and the test performs very well when $\kappa \geq 0.70$. This pattern is not particularly surprising given that the effective number of Fourier frequencies used to construct the MBLS estimator and sup-Wald test statistic over each subinterval is substantialler smaller when $\kappa = 0.50$ compared with $\kappa = 0.70$, as evidence by the corresponding "effective" full sample sizes of frequencies $|350^{0.50}| = 18$ and $|350^{0.70}| = 60$, respectively.

Generally, we find the LCM-based sup-Wald test, $\mathcal{P}_n(1)$, to perform very well across realistic persistence configurations and sample sizes. Moreover, the procedure is robust to changes tuning parameters, except when selecting the bandwidth parameter, κ , too low.

6 Empirical Analysis: Stability of the IV-RV Relation

As detailed in the introduction, and going back to Andersen & Bollerslev (1997), many empirical studies have documented the presence of long memory persistence in RV. Moreover, there is an extensive literature finding: (i) IV is an unbiased predictor of future RV after correcting for various methodological issues, and it subsumes the information content in historical RV, e.g., Christensen & Prabhala (1998) and Jiang & Tian (2005); (ii) IV and RV are fractionally cointegrated and their long-run relation is unbiased, e.g., Christensen & Nielsen (2006) and Bandi & Perron (2006); (iii) the fractionally cointegrating equilibrium between IV and RV carry information about the variance risk premium and forecast asset returns, e.g., Bollerslev et al. (2013) and Li et al. (2020).

In this section, we revisit the long-run relation between IV and RV by first examining the relative efficiency of the long-run information in IV for RV, comparing it to forecasts constructed from a longmemory model exploiting high-frequency data, thereby providing a more challenging benchmark than historical volatility, which the literature typically has relied upon. Second, we examine the stability of the fractional cointegrating relations using our new LCM-based sup-Wald tests. Finally, we explore the impact of "instability" on the forecasting prowess of the variance risk premium for returns.

6.1 Data and Summary Statistics

Our main dataset consists of three monthly series; implied volatility (IV) and realized volatility (RV) on the S&P 500 as well as a long-memory (LM) forecast of RV. We construct the RV series from oneminute observations of S&P 500 futures over January 1988 - December 2019. On each trading day, we use the flat-top realized kernel estimator of Varneskov (2016, 2017) to compute the intra-day quadratic variation, and we add the squared close-to-open return from the preceding trading day to account for the overnight return variation. We construct the RV series by summing all daily quadratic variation

estimates across the trading days in a given month, then apply a square-root transform, and annualize. The IV is proxied by the VIX. Finally, the LM forecasts are obtained using an $ARFIMA(1, d, 0)$ model. Specifically, we fit the model to daily log-transformed quadratic variation estimates and apply a lognormal correction to map the forecast back into levels. The latter is motivated by Andersen, Bollerslev, Diebold & Ebens (2001) and Andersen, Bollerslev, Diebold & Labys (2001, 2003), who show that RV is approximately log-normal and that long-memory models capture their dynamic properties well.⁷ The ARFIMA model is estimated using a two-year rolling window.⁸ The first two years of the sample are used to initialize the LM forecasts, so our final dataset spans January 1990 through December 2019, featuring $n = 360$ monthly observations. The series and their autocorrelation functions are depicted in Figure 1. Clearly, they are all highly persistent and strongly comoving, consistent with a large literature stipulating that IV and RV are fractionally cointegrated.

We further assess the properties of the series in Table 6, providing standard unconditional summary statistics and estimates of their degree of fractional integration. For the latter, we exploit the TELW estimator of Andersen & Varneskov (2020a) and the local polynomial Whittle with noise (LPWN) estimator by Frederiksen, Nielsen & Nielsen (2012), which is more robust against short-memory dynamics and measurement errors. The TELW estimator is implemented using a trimming parameter $\ell_d = \lfloor n^{0.30} \rfloor$ and a bandwidth $m_d = \lfloor n^{0.75} \rfloor$, whereas the LPWN estimator is implemented with a bandwidth $m_d = \lfloor n^{0.85} \rfloor$ and short-memory polynomials of orders zero and one. From Table 6, we observe that the unconditional mean of IV is larger than for RV, consistent with the presence of a volatility risk premium; see, e.g., Carr & Wu (2009) and Bollerslev et al. (2011). Second, RV is more volatile and has fatter tails than LM and IV, consistent with the former being a realization and the latter capturing volatility expectations. Finally, we confirm that the series are fractionally integrated, with orders falling in the range 0.5-0.6, when assessed through their LPWN estimates.

6.2 Econometric Issues and Implementation

There is an extensive literature exploring the IV–RV relation and correcting for various methodological issues. Although our exposition focuses on VIX, our adopted proxy for the risk-neutral expectation of IV, the following decomposition summarizes the econometric issues more broadly. We define the volatility risk premium for the coming month to be the difference between the expectation of RV under the risk-neutral measure, Q and the objective measure, P, yielding,

$$
\mathbf{VIX}_t = \mathbb{E}_t^{\mathbb{Q}}[\mathbf{RV}_{t+1}] + \mathcal{E}_t = \mathbb{E}_t^{\mathbb{P}}[\mathbf{RV}_{t+1}] + \underbrace{\mathbb{E}_t^{\mathbb{Q}}[\mathbf{RV}_{t+1}] - \mathbb{E}_t^{\mathbb{P}}[\mathbf{RV}_{t+1}]}_{\text{volatility risk premium}} + \mathcal{E}_t, \tag{26}
$$

where \mathcal{E}_t captures a Jensen's inequality bias from the square-root transformation and measurement errors in VIX, generated by a (time-varying) truncation of the strike range in the construction of the

⁷To corroborate this, we report unconditional summary statistics for both RV and log-RV in Table 6 below, which show that the logarithmic transformation alleviates skewness and kurtosis, rendering the series much closer to Gaussian. ⁸We have tested several different window lengths. The results are robust and, thus, omitted for brevity.

index and observation errors in the underlying options; see, e.g., Jiang & Tian (2005), Andersen & Bondarenko (2007), Andersen, Bondarenko & Gonzalez-Perez (2015) and Andersen, Fusari, Todorov & Varneskov (2020) .⁹ Hence, we have two sources of errors; one arising from standard measurement errors and another from the volatility risk premium. The latter is known to be positive on average (per our definition above), correlated with the level of volatility and weakly dependent; see, e.g., Bandi & Perron (2006), Chernov (2007), Bollerslev, Tauchen & Zhou (2009), Carr & Wu (2009) and Bollerslev et al. (2011). Next, we complement equation (26) with a decomposition of the latent P-measure forecast of RV,

$$
\mathbb{E}_t^{\mathbb{P}}\big[\text{RV}_{t+1}\big] \ \equiv \ \mathcal{V}_{t,1} \ + \ \mathcal{V}_{t,2} \,, \tag{27}
$$

where $V_{t,1} \in I(d_{\text{VIX}})$ and $V_{t,2} \in I(d_{\text{RES}})$ with $0 \leq d_{\text{RES}} < d_{\text{VIX}}$. That is, the latent RV forecast embedded in VIX has a two-factor structure, and the most persistent component is fractionally integrated, as documented in Table 6^{10} . The decompositions imply that VIX is fractionally integrated of order $I(d_{\text{VIX}})$, with the remaining components being of strictly smaller orders (often assumed to be $I(0)$). This motivated Bandi & Perron (2006), Christensen & Nielsen (2006), Nielsen (2007), Kellard et al. (2010), and Nielsen & Frederiksen (2011) to assess whether IV is "long-run unbiased" for RV, i.e., whether the most persistent forecast component, $V_{t,1}$, achieves a coefficient $\mathcal{B}_{IV} = 1$. These studies take advantage of the differing persistence among the VIX_t components and use NBLS-type estimators to identify the accuracy of the most persistent component. By the same logic, Bollerslev et al. (2011), Bollerslev et al. (2013), Osterrieder et al. (2019) and Li et al. (2020) specify fractional cointegration models between IV and RV to estimate the volatility risk premium component associated with this long-run equilibrium and use it to forecast asset returns.

While these studies, indeed, demonstrate that IV is long-run unbiased, they do not assess the efficiency of its predictions for the low-frequency movements in RV.¹¹ Our analysis goes one step further by examining the relative information embedded in the persistent IV component against a competing ARFIMA forecast constructed using high-frequency data, in addition to their associated biases. Moreover, we use our new testing framework to assess whether these long-run relations, exploring bias and efficiency, are invariant over time. To this end, we test for parameter (in)stability using our new LCMbased sup-Wald test, $\mathcal{P}_n(1)$. As explained above, the LCM procedure is robust to errors-in-variables (cf., Remark 1), as long as they are less persistent than the signal of interest. Hence, it facilitates robust inference and testing in this fractional cointegration setting.

We implement the LCM procedure using the LPWN estimator in the fractional filtering step with a bandwidth $\lfloor n^{0.85} \rfloor$, since it achieves a slightly faster convergence rate than TELW, and all the variables have $d < 1$. In addition, we follow existing semi-parametric studies of fractional cointegration and take the average integration order as a global d estimate. Table 6 verifies that this is innocuous, as

⁹The Jensen's inequality bias arises, because \texttt{VIX}_t^2 captures expectations about $\mathbb{E}_t^{\mathbb{Q}}[\texttt{RV}_{t+1}^2]$.

¹⁰Naturally, $\mathbb{E}_{t}^{\mathbb{P}}[\mathbf{RV}_{t+1}]$ may have more than two components, as long as $\mathcal{V}_{t,1}$ remains the more persistent one.

¹¹Among others, Christensen & Prabhala (1998), Jiang & Tian (2005) and Chernov (2007) find IV to subsume the information content in historical volatility for S&P 500, applying inference techniques resting on weakly dependent dynamics. Chraoenwong et al. (2009), Taylor et al. (2010) and Cheng & Fung (2012) extend these findings to other assets.

the individual estimates are close for LPWN. The MBLS estimator in the second step is configured with trimming and bandwidth rates $(\nu, \kappa) \in \{(0.2, 0.5), (0.2, 0.6), (0.3, 0.6), (0.3, 0.7)\}\,$ both for the full sample fractional cointegration analysis and for the LCM-based sup-Wald testing. Whereas the choice of ν is well-aligned with the recommendation in Andersen & Varneskov (2020a), the bandwidth rate is slightly smaller to emphasize robustness rather than efficiency, when estimating the cointegrating relation, by alleviating the impact of high-frequency components. These choices are in line with the values considered in the simulation study as well as the bounds provided by Table $1¹²$

In computing the asymptotic variance, we further set $\nu_G = \nu$ and $\kappa_G = \kappa + 0.2$, and use the TELW estimator to find b with corresponding trimming and bandwidth rates, $\nu_b = 0.30$ and $\kappa_b = 0.75$. The $\mathcal{P}_n(1)$ test is applied with $\epsilon = 0.2$, determining the search space Φ , to initialize the LCM estimates in the beginning and the end of the sample. The test results are, however, robust to this selection.

6.3 Implied and Realized Volatility

First, we carry out a full sample fractional cointegration analysis of the predictive relation between RV and either IV or LM. The regression results, assessing whether the two respective forecasting procedures are long-run unbiased, are reported in Table 7. The IV results, in particular, confirm the findings from the literature, demonstrating that its relation to RV is fractionally cointegrated with strength $0.4 \le b \le 0.5$. Moreover, we cannot reject that IV is long-run unbiased, $\mathcal{B}_{IV} = 1$. Note that the bounds on b imply convergence rates of $S_n(b) \asymp n^{\kappa/2 + \frac{2}{5}(1-\kappa)}$ and $S_n(b) \asymp \sqrt{n}$, respectively, as cointegration improves on the semi-parametric rate, \sqrt{m} , obtained for $b = 0$. The corresponding LCM results for the LM forecasts similarly demonstrate fractional cointegration with equivalent strength. However, these forecasts may be slightly upward biased at the lowest frequencies, as the Wald test rejects $\mathcal{B}_{LM} = 1$ at a 5% significance level, when $\kappa < 0.7$. This bias disappears for $\kappa = 0.7$.

Second, we provide full sample multivariate predictive regression results in Table 8. These show the LM forecasts emerging as more efficient than IV for capturing the low-frequency movements in RV. In particular, we reject $\mathcal{B}_{LM} = \mathcal{B}_{IV}$ at a 5% significance level for all combinations of tuning parameters and even fail to reject $(\mathcal{B}_{LM}, \mathcal{B}_{IV}) = (1, 0)$ in two of the four configurations. While these results appear to contradict Christensen & Prabhala (1998), Jiang & Tian (2005) and Chernov (2007), who find IV to be informationally efficient when assessed against historical volatility, it is important to note the subtle difference in interpretation. Our analysis only speaks to the relative quality of the lowfrequency signals in the variables. Hence, IV may have important high-frequency information, which is not captured by our cointegration analysis. Our empirical results, on the other hand, demonstrate that the most persistent component of IV may be well-approximated by an $ARFIMA(1, d, 0)$ process, generating more efficient predictions of the low-frequency variation in RV relative to IV.

Third, we test the stability of the predictive relations using the LCM-based sup-Wald test, with

¹²Whereas the lower bandwidth provides unbiased, yet less efficient, full sample LCM estimates, as seen in Andersen & Varneskov (2020a), our simulations show that the sup-Wald test will exhibit poor size properties when $\kappa = 0.50$. However, as the sup-Wald test size and the LCM coefficient estimate are both accurate when $\kappa = 0.70$, the use of multiple bandwidth parameters will allow us to assess robustness across different frequency ordinates.

results reported in Table 9. Importantly, the test overwhelmingly rejects the null hypothesis (4) for both univariate and multivariate predictions.¹³ These rejections imply that full sample coefficient estimates obtained from the LCM procedure in Tables 7-8 will reflect the average coefficients across regimes in the sample, as conveyed by Tables 11 and 15 in Appendix B (from the simulation study). Unfortunately, they will conceal any regime-specific information from the forecasting variables over the course of the sample. Hence, to explore the nature of the rejections in greater detail, we depict rolling window LCM coefficient estimates (with 95% confidence bands) in Figure 2. The window length is five years $(n = 60)$, and LCM is implemented with $(\nu, \kappa) = (0.3, 0.6)$. The figure illustrates parameter instability associated with the Asian crisis in 1997, the global financial crisis 2008-2009, and the low-volatility bull-market environment after 2013. In fact, from the rolling multivariate regressions we observe that both LM and IV carry important low-frequency information about RV during the periods from 1990 through 2004 and again from 2014 through 2019. However, the low-frequency information embedded in IV is completely subsumed in the period leading up to and including the financial crisis. These results have implications for volatility forecasting, indicating a shift in the low-frequency RV characteristics over the 30-year period, particularly during periods of financial turmoil.

Finally, these results have ramifications for asset pricing and return predictions. Specifically, most studies measure the volatility risk premium as $VRP_t = IV_t - RV_t$. Other studies estimate the corresponding cointegrating relation between IV_t and RV_t to more accurately recover the "long-run" component of the volatility risk premium; see, e.g., Bollerslev et al. (2011), Bollerslev et al. (2013), Osterrieder et al. (2019) and Li et al. (2020) .¹⁴ Our rejection of stability of the cointegration models for RV, thus, suggest that such measures may convey different information across predictive regimes. That is, since the equillibrium cointegrating relation and "long-run" risk premium may differ across the sample, the forecast variable VRP_t may contain different information across such regimes, being a mixture of long- and short-run components. As a simple gauge of this conjecture, we predict monthly returns for the S&P 500 using the VRP_t measure and standard OLS regressions.¹⁵ Specifically, we consider full sample predictions as well as subsamples from 1990.1-2004.12, 2005.1-2013.12 and 2014.1-2019.12, where IV either provides incremental low-frequency information about RV (first and third interval) or is subsumed by LM (second interval). Summary statistics for the return and VRP_t series are provided in Table 10 together with the predictive regressions results. Interestingly, we find that all predictive power from VRP_t materializes during the subsample 2005.1-2013.12, where the low-frequency information from IV is subsumed by time series forecasts. This suggests that it is the information from IV associated with high-frequency movements in RV, and thus the short-run component of the volatility

¹³Hence, despite the size distortions of the sup-Wald test $\mathcal{P}_n(1)$ for $\kappa = 0.50$, the results using the remaining tuning parameter configurations provide strongly significant evidence against stability of the cointegration parameter.

¹⁴Importantly, this does not change the definition of the volatility risk premium, $\mathbb{E}_t^{\mathbb{Q}}[\text{RV}_{t+1}] - \mathbb{E}_t^{\mathbb{P}}[\tilde{\text{RV}}_{t+1}]$. Indeed, the alternative ways of constructing the empirical forecast measures reflect an emphasis on identifying different components of the risk premium; namely, those associated with variation operating at different frequencies.

 15 Note that we implement the predictive return regressions assuming that both returns and the volatility risk premium are weakly dependent. A rigorous treatment of return regression using fractionally integrated forecast variables is given by Andersen & Varneskov (2020b). A related analysis of such issues is beyond the scope of this paper.

risk premium, which generates significant forecasting power for returns.

Hence, in addition to accommodating regime-dependent low-frequency dynamics in forecasting models for RV, our empirical results suggests that a potentially fruitful avenue for future research would be to extend existing asset pricing models with a volatility risk premium to allow for regimedependent cointegration between IV and RV as well as corresponding return predictability. Moreover, since, among others, Bollerslev & Todorov (2011), Andersen, Fusari & Todorov (2015) and Kilic & Shaliastovich (2019) have shown that the volatility risk premium is mainly associated with downside risk and/or left-tail jumps, it would be interesting to examine potential parameter instabilities in such persistent forecast variables as well. Again, this is left for future research.

7 Conclusion

This paper develops parameter instability and structural change tests for predictive regressions in economic systems governed by persistent vector autoregressive dynamics. Specifically, in a setting where all – or a subset – of the variables may be fractionally integrated, and the predictive relation may feature cointegration (possibly unknown, ex-ante), we provide sup-Wald break tests using the LCM approach of Andersen & Varneskov (2020a). The new tests cover both parameter variation and multiple structural changes with unknown break dates, and allows for the number of breaks to be unknown. The adoption of LCM ensures the tests, asymptotically, avoid problems associated with spurious inference, regression balance and endogenous variation in the predictors, which typically arise in such settings. We establish limit theory for the tests, showing it coincides with those for standard break testing procedures in, e.g., Andrews (1993) and Bai & Perron (1998), implying that existing critical values may be applied, without modification. Hence, the tests are easy to implement.

We apply the new structural change tests to explore the stability of the fractionally cointegrating relation between option implied and realized volatility. Moreover, we assess the relative efficiency of the low-frequency information embedded in IV forecasts against a challenging benchmark, a time series long-memory model-based forecast constructed from high-frequency data. Unlike existing studies, we find evidence that the IV-RV cointegrating relation is unstable, rejecting the null hypothesis of parameter stability at all conventional significance levels, and that carefully constructed time-series forecasts are more efficient than IV in capturing low-frequency movements in RV. Finally, we show that such parameter instabilities manifest themselves as instabilities in the forecasting relation between asset returns and the volatility risk premium.

	Tuning Parameter Bounds in Simulations										
		no cointegration: $b = 0$			cointegration: $b = d$						
	$d = 0.30$	$d = 0.60$	$d = 1.00$		$d = 0.30$	$d = 0.60$	$d = 1.00$				
ν_{\min}	0.240	0.180	0.125		$0.125 +$	$0.125 +$	$0.125 +$				
					$0.30(1 - \kappa)$	$0.60(1 - \kappa)$	$(1 - \kappa)$				
$\nu_{\rm max}$	κ	κ	κ		κ	κ	κ				
$\kappa_{\rm min}$	ν	ν	ν		$\nu \vee 0.375$	$\nu \vee 6/11$	$\nu \vee 2/3$				
$\kappa_{\rm max}$	4/5	4/5	4/5		4/5	4/5	4/5				

Table 1: Tuning parameter bounds. This table compiles the lower and upper bounds on the trimming and bandwidth rates required by Assumption T for the different configurations of d and b considered for the simulation study in Section 5. Specifically, the bounds are derived using $\rho = 0.75$ for the TELW estimator. Moreover, for simplicity of exposition, the lower trimming rate bound, ν_{\min} , is derived under the assumption $\kappa \geq 1/2$ to avoid the appearance of multiple terms, as in equation (12). The lower bound ν_{\min} for $d \leq 0.60$ and $b = 0$ follows from the first restriction in equation (12), that is, from $(3/5 - (2d - b)/5)/2 < \nu$, and the remaining cases from $(1 - \varrho)/2 + b(1 - \kappa) < \nu$.

			Size and Power of the LCM-based Sup-Wald Test: Baseline						
				no cointegration: $b = 0$					
		\mathcal{H}_0 , $n=$		$\mathcal{H}_1, n =$	\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$	
	350	700	350	700	350	700		350	700
$d = 0.30$	0.0850	0.0690	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000
$d = 0.60$	0.0850	0.0690	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000
$d = 1.00$	0.0850	0.0690	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000
				cointegration: $b = d$					
		\mathcal{H}_0 , $n=$		$\mathcal{H}_1, n =$	\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$	
	350	700	350	700	350	700		350	700
$d = 0.30$	0.0610	0.0640	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000
$d = 0.60$	0.0610	0.0470	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000
$d = 1.00$	0.0260	0.0200	1.0000	1.0000	1.0000	1.0000		1.0000	1.0000

Table 2: Size and power: Baseline. This table provides size and power results for the LCM-based sup-Wald test (19), which is robust against general forms of parameter instability. Specifically, the test is implemented using trimming $\ell = \lfloor n^{\nu} \rfloor$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. Robustness towards the tuning parameters is studied in Table 4. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_0 describes the null hypothesis *without* parameter instability, i.e., it speaks to test size. \mathcal{H}_1 , \mathcal{H}_2 , \mathcal{H}_3 describes three baseline predictive models with breaks, thus illustrating the power of the test. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

			Size and Power of the LCM-based Sup-Wald Test: Small Breaks					
					no cointegration: $b = 0$			
		\mathcal{H}_0 , $n=$		\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$		\mathcal{H}_6 , $n=$
	350	700	350	700	350	700	350	700
$d = 0.30$	0.0850	0.0690	0.4740	0.7010	0.5840	0.8110	0.4830	0.7160
$d = 0.60$	0.0850	0.0690	0.4740	0.7010	0.5840	0.8110	0.4830	0.7160
$d = 1.00$	0.0850	0.0690	0.4740	0.7010	0.5840	0.8110	0.4830	0.7160
					cointegration: $b = d$			
		\mathcal{H}_0 , $n=$		\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$		\mathcal{H}_6 , $n=$
	350	700	350	700	350	700	350	700
$d = 0.30$	0.0610	0.0640	0.4870	0.8120	0.6060	0.8770	0.4310	0.7420
$d = 0.60$	0.0610	0.0470	0.5470	0.8850	0.6320	0.9480	0.4310	0.7930
$d = 1.00$	0.0260	0.0200	0.4860	0.8960	0.5630	0.9500	0.3370	0.7440

Table 3: Size and power: Small breaks. This table provides size and power results for the LCM-based sup-Wald test (19), which is robust against general forms of parameter instability. Specifically, the test is implemented using trimming $\ell = \lfloor n^{\nu} \rfloor$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. Robustness towards the tuning parameters is studied in Table 5. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_0 describes the null hypothesis *without* parameter instability, i.e., it speaks to test size. H_4 , H_5 , H_6 describes three predictive models with small breaks, thus illustrating the power of the test. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

Table 4: Size and power: Tuning parameters, baseline. This table provides size and power results for the LCM-based sup-Wald test (19), which is robust against general forms of parameter instability. Specifically, the test is implemented using several combinations of the tuning parameters; trimming $\ell = \lfloor n^{\nu} \rfloor$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . H₀ describes the null hypothesis *without* parameter instability, i.e., it provides size results. \mathcal{H}_1 , \mathcal{H}_2 , \mathcal{H}_3 describes three baseline predictive models with breaks, thus speaking to the power of the test. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is $n = 350$, the integration order is $d = 0.60$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

	Size and Power of the LCM-based Sup-Wald Test: Tuning Parameters, Small Breaks											
		no cointegration: $b = 0$				cointegration: $b = d$						
$(\nu,\kappa,\kappa_G) =$	\mathcal{H}_0	\mathcal{H}_4	\mathcal{H}_5	\mathcal{H}_6		\mathcal{H}_0	\mathcal{H}_4	\mathcal{H}_5	\mathcal{H}_6			
(0.10, 0.70, 0.90)	0.0710	0.4710	0.5780	0.4720		0.0530	0.5360	0.6330	0.4210			
(0.20, 0.70, 0.90)	0.0850	0.4740	0.5840	0.4830		0.0610	0.5470	0.6320	0.4310			
(0.30, 0.70, 0.90)	0.0890	0.4890	0.5950	0.4970		0.0750	0.5450	0.6420	0.4470			
(0.20, 0.50, 0.90)	0.1580	0.3350	0.4150	0.3610		0.2020	0.6530	0.7080	0.5060			
(0.20, 0.60, 0.90)	0.1090	0.3750	0.4550	0.3780		0.0970	0.5800	0.6680	0.4620			
(0.20, 0.75, 0.90)	0.0800	0.5650	0.6870	0.5670		0.0450	0.5560	0.6590	0.4330			
(0.20, 0.70, 0.75)	0.0800	0.4810	0.5880	0.4840		0.0710	0.6280	0.7200	0.5190			
(0.20, 0.70, 0.80)	0.0750	0.4770	0.5860	0.4720		0.0580	0.5850	0.6880	0.4800			

Table 5: Size and power: Tuning parameters, small breaks. This table provides size and power results for the LCM-based sup-Wald test (19), which is robust against general forms of parameter instability. Specifically, the test is implemented using several combinations of the tuning parameters; trimming $\ell = \lfloor n^{\nu} \rfloor$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . H₀ describes the null hypothesis *without* parameter instability, i.e., it speaks to the test size. \mathcal{H}_4 , \mathcal{H}_5 , \mathcal{H}_6 describes three predictive models with *small* breaks, thus illustrating the power of the test. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is $n = 350$, the integration order is $d = 0.60$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

Figure 1: Volatility Series and Autocorrelation Functions (ACFs.) This figure depicts the series and the ACFs for the realized volatility (RV) series as well as the ARFIMA (LM) and model-free implied volatility (IV) forecasts. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$). The sample autocorrelation functions are computed for the first 100 lags for each variable in the full sample.

	Summary Statistics for Volatility Measures											
	mean	StDev	Skewness	Kurtosis	TELW	LPWN0	LPWN1					
RV	0.1509	0.0826	2.9141	18.0456	0.5566 (0.0552)	0.4763 (0.0809)	0.5098 (0.1217)					
LM	0.1403	0.0667	1.9758	9.6354	0.6309 (0.0552)	0.6167 (0.0751)	0.6101 (0.1118)					
ΙV	0.1919	0.0737	1.7536	7.7572	0.6957 (0.0552)	0.5904 (0.0720)	0.5423 (0.1181)					
$log-RV$	-1.9978	0.4392	0.6457	3.6146	0.5297 (0.0552)	0.5478 (0.0766)	0.5731 (0.1150)					

Table 6: Summary statistics. This table displays summary statistics for the realized volatility (RV) series as well as the ARFIMA (LM) and model-free implied volatility (IV) forecasts. Furthermore, we include corresponding summary statistics for log-RV, since the monthly LM forecast are generated from daily log-transformed RV estimates. In addition to the standard unconditional statistics, we present estimates of the fractional integration order for all variables using the trimmed exact local Whittle (TELW) estimator from Andersen & Varneskov (2020a) and the local polynomial Whittle with noise (LPWN) estimator from Frederiksen et al. (2012), with standard errors indicated in parentheses. The TELW estimator is implemented using trimming $\ell_d = \lfloor n^{0.30} \rfloor$ and bandwidth $m_d = \lfloor n^{0.75} \rfloor$. The LPWN estimator is implemented with bandwidth $m_d = \lfloor n^{0.85} \rfloor$. The respective numbers for the LPWN signifies the order of the short-memory polynomial. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$).

Univariate LCM Cointegration Analysis										
			LM, $(\nu, \kappa) =$							
	(0.2, 0.5)	(0.2, 0.6)	(0.3, 0.6)	(0.3, 0.7)	(0.2, 0.5)	(0.2, 0.6)	(0.3, 0.6)	(0.3, 0.7)		
\mathcal{B}	1.1717 (15.0419)	1.1664 (14.7894)	1.1736 (14.8583)	1.0616 (12.5238)	0.9638 (11.8750)	0.9421 (11.2118)	0.9540 (11.3776)	0.8787 (9.6458)		
b	0.4906 (8.8855)	0.4898 (8.8855)	0.4909 (8.8909)	0.4620 (8.3677)	0.4478 (8.1099)	0.4388 (7.9462)	0.4438 (8.0368)	0.4106 (7.4363)		
WALD	4.8587	4.4523	4.8322	0.5288	0.1993	0.4750	0.3010	1.7738		
p -WALD	0.0275	0.0349	0.0279	0.4671	0.6552	0.4907	0.5832	0.1829		

Table 7: Univariate LCM analysis. This table displays results from a univariate LCM analysis, where RV is predicted using either ARFIMA (LM) or model-free implied volatility (IV) forecasts. Specifically, the coefficient estimate and associated t-statistic (in parenthesis) is provided along with a TELW estimate of the fractional cointegration strength, b, and associated t-statistics (in parenthesis) and, finally, a Wald test of $\mathcal{B} = 1$ together with its p-values. The results are provided for $\hat{\mathcal{B}}(\ell, m)$, where several combinations of trimming $\ell = \lfloor n^{\nu} \rfloor$ and bandwidth $m = |n^{\kappa}|$ parameters are considered. The TELW estimator is implemented as described in Table 6. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$).

	Multivariate LCM Cointegration Analysis												
$(\nu,\kappa) =$	\mathcal{B}_{LM}	$\mathcal{B}_{\rm IV}$	b	W1	$p-W1$	W2	$p-W2$	W3	p -W3				
(0.2, 0.5)	1.0627 (6.1677)	0.0969 (0.6130)	0.5077 (9.1946)	4.6694	0.0968	38.198	0.0000	11.783	0.0006				
(0.2, 0.6)	0.8389 (6.4918)	0.2905 (2.2244)	0.5097 (9.2310)	5.8705	0.0531	42.219	0.0000	6.3382	0.0118				
(0.3, 0.6)	0.8170 (6.3337)	0.3198 (2.4524)	0.5125 (9.2812)	6.9762	0.0306	40.132	0.0000	5.2258	0.0223				
(0.3, 0.7)	0.9232 (8.1353)	0.1326 (0.9696)	0.4813 (8.7168)	0.9401	0.6250	67.009	0.0000	14.732	0.0001				

Table 8: Multivariate LCM analysis. This table displays results from a multivariate LCM analysis, where RV is predicted using ARFIMA (LM) and model-free implied volatility (IV) forecasts. The coefficient estimates and t-statistics (in parentheses) are provided along with a TELW estimate of the fractional cointegration strength, b, and associated t-statistics (in parenthesis). Finally, the Wald test statistics and p -values from the three hypotheses $(\mathcal{B}_{LM}, \mathcal{B}_{IV}) = (1, 0), (\mathcal{B}_{LM}, \mathcal{B}_{IV}) = (0, 1)$ and $\mathcal{B}_{LM} = \mathcal{B}_{IV}$ are reported, labeled W1, W2 and W3 as well as p-W1, p-W2 and p-W3, respectively. The results are provided for $\hat{\mathcal{B}}(\ell, m)$, where several combinations of trimming $\ell = \lfloor n^{\nu} \rfloor$ and bandwidth $m = \lfloor n^{\kappa} \rfloor$ parameters are considered. The TELW estimator is implemented as in Table 6. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$).

Testing Stability of IV-RV Relations											
	Sup-Wald test, $(\nu, \kappa) =$										
	(0.3, 0.7) (0.2, 0.5) (0.2, 0.6) (0.3, 0.6)										
LM	$16.5755***$	19.3363***	$19.6658***$	$15.8335***$							
IV	$9.4495**$	$10.7750**$	13.7734***	11.3375**							
$LM + IV$	25.7693***	$41.4712***$	37.5787***	$28.1861***$							

Table 9: Stability testing. This table displays results of the sup-Wald test for stability of univariate and multivariate LCM cointegrating regressions for RV. The latter is predicted using either ARFIMA (LM) and/or model-free implied volatility (IV) forecasts. Specifically, the sup-Wald test is implemented using $\epsilon = 0.2$. The critical values of the test may be found in Andrews (1993, Table 1) for both univariate and multivariate regressions, with degrees of freedoms equal to $k = 1$ and $k = 2$, respectively. The results are provided for $\mathcal{B}(\ell, m)$, where several combinations of trimming $\ell = |n^{\nu}|$ and bandwidth $m = |n^{\kappa}|$ parameters are considered. The TELW estimator is implemented as described in Table 6, and the variables are fractionally filtered using the full-sample global estimate of the integration order prior to sample splits and MBLS estimation. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$). Finally, ** and *** signify rejection of the null hypothesis, \mathcal{H}_0 , of parameter stability at the 5% and 1% significance level, respectively.

Figure 2: LCM parameter estimates. This figure depicts rolling window LCM estimates and 95% confidence bands from regressions of realized volatility (RV) series on ARFIMA (LM) and model-free implied volatility (IV) forecasts. All variables are in square-root form. The sample spans January 1990 through December 2019 ($n = 360$). The window length is 5 years ($n = 60$), and LCM is implemented with (ν, κ) = (0.3, 0.6).

Return Predictions using VRP										
Panel A: Summary	mean	StDev	Skewness	Kurtosis						
returns	0.0063	0.0412	-0.8109	4.7860						
VRP	0.0407	0.0376	-1.5217	11.0565						
Panel B: Regressions		\mathcal{T}_1	\mathcal{T}_2	\mathcal{T}_3						
constant	-0.0003 (-0.0945)	0.0051 (1.1293)	-0.0093 (-1.8588)	0.0083 (1.5307)						
B_{VRP}	0.1641 (2.5570)	0.0478 (0.6217)	0.3430 (4.9452)	-0.0191 (-0.1656)						
adj. \mathbb{R}^2	0.0170	-0.0098	0.1009	-0.0287						

Table 10: Return regressions. This table displays summary statistics for monthly $S\&P$ 500 returns and volatility risk premium (VRP) estimates in Panel A along with standard predictive return regression results in Panel B. Specifically, the latter is based on OLS estimates and the heteroskedasticity and autocorrelation consistent (HAC) inference procedure by Andrews (1991) using the Parzel kernel and a bandwidth $\frac{4(n/100)^{2/9}}{1}$. The summary statistics and regressions results for " T " are based on the full sample from January 1990 through December 2019 $(n = 360)$. Regression results are also provided for three subsamples from 1990.1-2004.12, 2005.1-2013.12 and 2014.1-2019.12, labeled \mathcal{T}_1 , \mathcal{T}_2 and \mathcal{T}_3 , respectively. These subsamples correspond approximately to the periods in Figure 2, where the multivariate coefficient estimate on IV is positive, turns negative, and then back to positive.

A Proofs

This section contains the proofs of the asymptotic results in the paper. Before proceeding, we introduce some notation. For a generic vector V , let $V(i)$ index the *i*th element, and, similarly, for a matrix M, let $M(i, q)$ denote its (i, q) th element. Moreover, denote by $K \in (0, \infty)$ a generic constant, which may take different values from line to line or from (in)equality to (in)equality. Finally, we remark that sometimes the (stochastic) orders refer to scalars, sometimes to vectors and matrices.

A.1 Proof of Theorem 1

The proof relies heavily on asymptotic results and stochastic bounds from the corresponding proof of the full sample LCM result in the Online Appendix to Andersen & Varneskov (2020a). We will henceforth refer to the latter as AVOA (2020). First, by applying the decomposition in AVOA (2020, equation $(A.2)$) in conjunction with AVOA (2020, Lemma $A.2(b)$), we may write,

$$
\sqrt{m_g} \lambda_{m_g}^{-b}(g) \left\| \widehat{\mathcal{B}}(\ell,m,g) - \mathcal{B} - \widehat{F}_{\widehat{u}\widehat{u}}(\ell,m,g)^{-1} \widehat{F}_{u,\eta}^{(t,1,b)}(1,m,g) \right\| \xrightarrow{\mathbb{P}} 0, \tag{A.1}
$$

for all partitions $g = 0, \ldots, q$, where $\widehat{F}_{u,\eta}^{(t,1,b)}(1,m,g)$ is the TDAC between u_{t-1} and $\eta_t^{(b)}$ $t^{(0)}$. Moreover, by applying AVOA $(2020, \text{Lemma A}.2(a))$, we have,

$$
\left\|\lambda_{m_g}^{-1}(g)\widehat{F}_{\widehat{u}\widehat{u}}(\ell,m,g)-\boldsymbol{G}_{uu}\right\| \stackrel{\mathbb{P}}{\to} 0,
$$
\n(A.2)

similarly, for all $g = 0, \ldots, q$, where G_{uu} is positive definite by Assumption D1. Next, define

$$
\widetilde{F}_{u,\eta}^{(t,1,b)}(1,m,g) \equiv \frac{2\pi}{n_g} \sum_{j=1}^{m_g} \Re\left(\widetilde{I}_{u\eta}^{(t,1,b)}(\lambda_j,g)\right), \quad \widetilde{I}_{u\eta}^{(t,1,b)}(\lambda_j,g) = \boldsymbol{w}_u(\lambda_j,g)\lambda_j^b(g)e^{(\pi/2)bi}\bar{w}_\eta(\lambda_j,g), \quad \text{(A.3)}
$$

where $w_u(\lambda_j, g)$ and $w_\eta(\lambda_j, g)$ are the discrete Fourier transforms of u_{t-1} and $\eta_t^{(b)}$ $t^{(0)}$, respectively. Moreover, define some arbitrary vector $\boldsymbol{\psi} = (\psi_1, \dots, \psi_k)'$ as well as,

$$
\mathcal{U}_4(g,b) \equiv \sum_{t \in \mathcal{T}_g} \epsilon'_t \sum_{s=1}^{t-1} C^n_{t-s}(g,b) \epsilon_s, \quad C^n_t(g,b) \equiv \frac{1}{2\pi n_g \sqrt{m_g}} \sum_{j=1}^{m_g} \lambda_j^b(g) \chi_j(g,b), \tag{A.4}
$$

where the sequence of coefficients, $\chi_j(g, b)$, is defined as $\chi_j(g, b) \equiv \sum_{i=1}^k \lambda_{m_g}^{-b}(g) \psi_i \mathcal{A}_j(g, b, i)$, with

$$
\mathcal{A}_{j}(g,b,i) \equiv \Re \left(\left(\mathbf{A}(\lambda_{j}(g),i)^{\prime} \bar{\mathbf{A}}(\lambda_{j}(g),k+1)^{\prime} e^{-i(t-s)\lambda_{j}(g)} \right) e^{(\pi/2)bi} \right) + \Re \left(\left(\mathbf{A}(\lambda_{j}(g),k+1)^{\prime} \bar{\mathbf{A}}(\lambda_{j}(g),i)^{\prime} e^{i(t-s)\lambda_{j}(g)} \right) e^{(\pi/2)bi} \right).
$$

Then, by applying AVOA (2020, Lemma A.11), we have,

$$
\sqrt{m_g} \lambda_{m_g}^{-b}(g) \left\| \widehat{F}_{u,\eta}^{(t,1,b)}(1,m,g) - \widetilde{F}_{u,\eta}^{(t,1,b)}(1,m,g) \right\| \xrightarrow{\mathbb{P}} 0, \tag{A.5}
$$

for all partitions $g = 0, \ldots, q$. Moreover, by using the decomposition in AVOA (2020, Equation (A.16)) in conjunction with the same arguments for their terms $\mathcal{U}_1, \mathcal{U}_2, \mathcal{U}_3$ and \mathcal{U}_4 , it follows,

$$
\left|\sqrt{m_g}\lambda_{m_g}^{-b}(g)\psi'\widetilde{F}_{u,\eta}^{(t,1,b)}(1,m,g)-\mathcal{U}_4(g,b)\right| \xrightarrow{\mathbb{P}} 0, \tag{A.6}
$$

similarly, for all $g = 0, \ldots, q$. Hence, since $\mathcal{M}_t(g, b) \equiv \epsilon'_t \sum_{s=1}^{t-1} C_{t-s}^n(g, b) \epsilon_s$ in $\mathcal{U}_4(g, b)$ is a martingale difference sequence with respect to the filtration \mathcal{F}_{t-1} , we use the same arguments as for AVOA (2020, Lemma A.3) to show,

$$
\mathcal{U}_4(g,b) \xrightarrow{\mathbb{D}} N\left(0, \boldsymbol{\psi}^{\prime} \boldsymbol{G}_{uu} \boldsymbol{\psi} \frac{G_{\eta\eta}}{2(1+2b)}\right), \tag{A.7}
$$

which applies, again, for all $g = 0, \ldots, q$. Next, as $\mathcal{M}_t(g, b)$ is a martingale difference sequence and the sample partitions \mathcal{T}_g , $g = 0, \ldots, q$, are non-overlapping, it readily follows that the covariance between $\mathcal{U}_4(g, b)$ and $\mathcal{U}_4(h, b)$ is trivial, when $g \neq h$. Hence, we may apply $(A.1)-(A.7)$ in conjunction with the continuous mapping theorem, Slutsky's theorem and the Cram´er-Wold theorem to show that,

$$
\sqrt{m}\lambda_m^{-b}\begin{pmatrix}\widehat{\mathcal{B}}(\ell,m,0)-\mathcal{B} \\
\vdots \\
\widehat{\mathcal{B}}(\ell,m,q)-\mathcal{B}\end{pmatrix}\xrightarrow{\mathbb{D}} N\left(\mathbf{0},\widetilde{\boldsymbol{\vartheta}}(b)\circ\mathbf{C}(b)\right), \quad \text{where}
$$
\n(A.8)

$$
\widetilde{\boldsymbol{\vartheta}}(b) = \text{diag}\left[(\varphi_1 - \varphi_0)^{2b(\kappa - 1) - \kappa}, \dots, (\varphi_{q+1} - \varphi_q)^{2b(\kappa - 1) - \kappa} \right],\tag{A.9}
$$

holds finite-dimensionally. Next, we establish *tightness* in the sample partition φ of the functional limit sequence. To this end, let us define $\widetilde{\mathcal{U}}_4(g, b) \equiv (\sqrt{m/m_g} \lambda_{m_g}^b \lambda_{m_g}^{-b}) \mathcal{U}_4(g, b)$. Then, since it follows that $\sqrt{m/m_g} \lambda_{m_g}^b \lambda_m^{-b} \to (\varphi_{g+1} - \varphi_g)^{b(\kappa-1)-\kappa/2} > 0$ for all partitions $g = 0, \ldots, q$ on the set Φ , we have that the convergence results in $(A.1)$, $(A.2)$, $(A.5)$ and $(A.6)$ still apply when taking the supremum over the partition vector φ on the set Φ . Hence, we may study tightness of the joint limit theory via the re-scaled variable $\mathcal{U}_4(g, b)$. By the law of iterated expectations and conditional independence of the martingale difference sequences over disjoint sets of observations, we have

$$
\mathbb{E}\left[\left(\widetilde{\mathcal{U}}_4(g,b)-\widetilde{\mathcal{U}}_4(h,b)\right)^2\right] \leq K\left((\varphi_{g+1}-\varphi_g)^{b(\kappa-1)-\kappa/2}-(\varphi_{h+1}-\varphi_h)^{b(\kappa-1)-\kappa/2}\right)^2
$$

$$
\leq K((\varphi_{g+1}-\varphi_g)-(\varphi_{h+1}-\varphi_h))^{2b(\kappa-1)-\kappa}, \tag{A.10}
$$

where the second inequality follows from the algebraic relation $|a + b|^p - |a|^p \le |b|^p$ for $p \in (0, 1)$, along with addition and subtraction. Hence, by applying (A.10) together with (A.8) and Ibragimov & Hasminskii (1981, Theorem 20), the sequence is tight and the joint uniform limit theory holds, whose representation in terms of independent Brownian motions follows by Wooldridge & White (1988,

Corollary 4.2), with their conditions D.1 and D.2 being satisfied by $(A.8)$, tightness, and Assumptions D1-D3. Finally, the mutual consistency condition follows exactly as in AVOA (2020, p. 2). \Box

A.2 Proof of Theorem 2

Part (a) follows by applying Assumption B and AVOA (2020, Lemma A.4) in conjunction with the continuous mapping theorem for all partition $g = 0, \ldots, q$ of every sequence of φ on Φ . (b) follows by applying (a), Theorem 1, the continuous mapping theorem and Slutsky's theorem. \Box

A.3 Proof of Theorem 3

First, by applying Theorem 1, we may write,

$$
S_n(b) \left(\hat{\beta}(\ell, m, 0) - \hat{\beta}(\ell, m, 1) \right) \Rightarrow C_{k_{\beta}}(b)^{1/2} \left(\frac{\mathbf{Y}(\varphi_1)}{\varphi_1^{\zeta + 1/2}} - \frac{(\mathbf{Y}(1) - \mathbf{Y}(\varphi_1))}{(1 - \varphi_1)^{\zeta + 1/2}} \right)
$$

=
$$
C_{k_{\beta}}(b)^{1/2} \left(\frac{\mathbf{Y}(\varphi_1)(1 - \varphi_1)^{\zeta + 1/2} - (\mathbf{Y}(1) - \mathbf{Y}(\varphi_1))\varphi_1^{\zeta + 1/2}}{\varphi_1^{\zeta + 1/2}(1 - \varphi_1)^{\zeta + 1/2}} \right)
$$
(A.11)

with, again, $S_n(b) = \sqrt{m} \lambda_m^{-b}$, and where $C_{k_\beta}(b)$ is the upper-left $k_\beta \times k_\beta$ submatrix of $C(b)$. Next, by applying Theorem 2 and the continuous mapping theorem,

$$
S_n(b)^2 \left(\widehat{\text{AVAR}}_{\beta}(0, b) + \widehat{\text{AVAR}}_{\beta}(1, b) \right)
$$

$$
\xrightarrow{\mathbb{P}} \left(\frac{1}{\varphi_1^{2\zeta}} + \frac{1}{(1 - \varphi_1)^{2\zeta}} \right) \mathbf{C}_{k_{\beta}}(b) = \left(\frac{(1 - \varphi_1)^{2\zeta} + \varphi_1^{2\zeta}}{\varphi_1^{2\zeta} (1 - \varphi_1)^{2\zeta}} \right) \mathbf{C}_{k_{\beta}}(b).
$$
 (A.12)

Hence, the final convergence results follow by combining $(A.11)$ and $(A.12)$ with the continuous mapping theorem and Slutsky's theorem, for both statements. \Box

A.4 Proof of Theorem 4

The result follows by applying the same arguments used to establish Theorem 3 for each partition $g = 0, \ldots, q$ on Φ in conjunction with asymptotic independence between partitions, by Theorem 1. \Box

B Additional Simulation Results

Tables 11-14 provide supplementary results for H_0 and the baseline alternatives H_1 , H_2 and H_3 . Specifically, they display the mean and standard deviation of the "full sample" LCM coefficient estimates to gauge the impact of breaks under the alternative hypotheses as well as the bias and RMSE for the estimated break fraction, φ_1 , using the LCM-based sup-Wald procedure. Tables 15-18 contain corresponding results for the small break alternatives.

			Mean LCM Estimate: Baseline						
				No Cointegration: $b = 0$					
		\mathcal{H}_0 , $n=$		\mathcal{H}_1 , $n=$			\mathcal{H}_2 , $n=$	\mathcal{H}_3 , $n=$	
	350	700	350	700		350	700	350	700
$d = 0.30$	0.9978	0.9983	1.2493	1.2477		0.9955	0.9966	0.7469	0.7463
$d = 0.60$	0.9978	0.9983	1.2493	1.2477		0.9955	0.9966	0.7469	0.7463
$d = 1.00$	0.9978	0.9983	1.2493	1.2477		0.9955	0.9966	0.7469	0.7463
				Cointegration: $b = d$					
		\mathcal{H}_0 , $n=$		$\mathcal{H}_1, n =$		\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$
	350	700	350	700		350	700	350	700
$d = 0.30$	0.9984	0.9992	1.2498	1.2486		0.9960	0.9974	0.7473	0.7470
$d = 0.60$	0.9987	0.9997	1.2500	1.2491		0.9962	0.9979	0.7475	0.7475
$d = 1.00$	0.9990	1.0002	1.2503	1.2496		0.9965	0.9983	0.7477	0.7479

Table 11: Mean estimate: Baseline. This table reports the mean of the full-sample LCM estimates for different predictive models. Specifically, the estimator is implemented using trimming $\ell = |n^{\nu}|$ and bandwidth $m = |n^{\kappa}|$, for $(\nu, \kappa) = (0.2, 0.7)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_0 describes the null hypothesis without parameter instability. H_1 , H_2 , H_3 describes three different baseline predictive models with breaks. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

			Standard deviation of LCM Estimate: Baseline								
				No Cointegration: $b=0$							
		$\mathcal{H}_0, n =$	$\mathcal{H}_1, n =$				\mathcal{H}_2 , $n=$		\mathcal{H}_3 , $n=$		
	350	700	350	700	350		700		350	700	
$d = 0.30$	0.0471	0.0366	0.0730	0.0562	0.0802		0.0603		0.0727	0.0545	
$d = 0.60$	0.0471	0.0366	0.0730	0.0562	0.0802		0.0603		0.0727	0.0545	
$d = 1.00$	0.0471	0.0366	0.0730	0.0562	0.0802		0.0603		0.0727	0.0545	
				Cointegration: $b = d$							
		\mathcal{H}_0 , $n=$		\mathcal{H}_1 , $n=$		\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$		
	350	700	350	700	350		700		350	700	
$d = 0.30$	0.0388	0.0295	0.0678	0.0521	0.0760		0.0566		0.0673	0.0502	
$d = 0.60$	0.0335	0.0241	0.0648	0.0494	0.0738		0.0545		0.0643	0.0476	
$d = 1.00$	0.0291	0.0185	0.0625	0.0469	0.0722		0.0530		0.0622	0.0456	

Table 12: Standard deviation: Baseline. This table provides the standard deviation of the full sample LCM estimates for different predictive models. Specifically, the estimator is implemented using trimming $\ell = \lfloor n^{\nu} \rfloor$ and bandwidth $m = \lfloor n^{\kappa} \rfloor$, for $(\nu, \kappa) = (0.2, 0.7)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_0 describes the null hypothesis without parameter instability. \mathcal{H}_1 , \mathcal{H}_2 , \mathcal{H}_3 describes three different baseline predictive models with breaks. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

						Bias of LCM-based Break Fraction Estimate: Baseline						
		No Cointegration: $b = 0$										
	\mathcal{H}_1 , $n=$				\mathcal{H}_2 , $n=$	\mathcal{H}_3 , $n=$						
	350 700			700 350		350	700					
$d = 0.30$	-0.0060	-0.0031		0.0002	-0.0001	0.0023	0.0031					
$d = 0.60$	-0.0060	-0.0031		0.0002	-0.0001	0.0023	0.0031					
$d = 1.00$	-0.0060	-0.0031		0.0002	-0.0001	0.0023	0.0031					
					Cointegration: $b = d$							
		\mathcal{H}_1 , $n=$			\mathcal{H}_2 , $n=$	\mathcal{H}_3 , $n=$						
	350	700		350	700	350	700					
$d = 0.30$	-0.0042	-0.0014		-0.0008	-0.0005	-0.0012	0.0003					
$d = 0.60$	-0.0031	-0.0010		-0.0021	-0.0026	-0.0027	-0.0013					
$d = 1.00$	-0.0033	-0.0016		-0.0032	-0.0042	-0.0046	-0.0016					

Table 13: Break fraction bias: Baseline. This table reports the bias of the break fraction estimates from the LCM-based sup-Wald test in equation (19). Specifically, the test is implemented using trimming $\ell = |n^{\nu}|$, bandwidth $m = \lfloor n^{\kappa} \rfloor$, and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_1 , \mathcal{H}_2 , \mathcal{H}_3 describes three different baseline predictive models with breaks, whose break fractions are 0.25, 0.50 and 0.75, respectively. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

RMSE of LCM-based Break Fraction Estimate: Baseline											
	No Cointegration: $b = 0$										
	\mathcal{H}_1 , $n=$			\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$				
	350	700		350	700	350	700				
$d = 0.30$	0.0160	0.0107		0.0141	0.0088	0.0162	0.0113				
$d = 0.60$	0.0160	0.0107		0.0141	0.0088	0.0162	0.0113				
$d = 1.00$	0.0160	0.0107		0.0141	0.0088	0.0162	0.0113				
	Cointegration: $b = d$										
	\mathcal{H}_1 , $n=$			\mathcal{H}_2 , $n=$			\mathcal{H}_3 , $n=$				
	350	700		350	700	350	700				
$d = 0.30$	0.0161	0.0116		0.0159	0.0099	0.0186	0.0120				
$d = 0.60$	0.0147	0.0092		0.0152	0.0099	0.0190	0.0120				
$d = 1.00$	0.0133	0.0080		0.0155	0.0122	0.0216	0.0128				

Table 14: Break fraction RMSE: Baseline. This table reports the bias of the break fraction estimates from the LCM-based sup-Wald test in equation (19). Specifically, the test is implemented using trimming $\ell = |n^{\nu}|$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_1 , \mathcal{H}_2 , \mathcal{H}_3 describes three different baseline predictive models with breaks, whose break fractions are 0.25, 0.50 and 0.75, respectively. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

Mean LCM Estimate: Small Breaks												
	No Cointegration: $b = 0$											
	\mathcal{H}_0 , $n=$			\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$			
	350	700	350	700		350	700		350	700		
$d = 0.30$	0.9978	0.9983	1.0482	1.0482		0.9976	0.9981		0.9479	0.9480		
$d = 0.60$	0.9978	0.9983	1.0482	1.0482		0.9976	0.9981		0.9479	0.9480		
$d = 1.00$	0.9978	0.9983	1.0482	1.0482		0.9976	0.9981		0.9479	0.9480		
	Cointegration: $b = d$											
	\mathcal{H}_0 , $n=$			\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$			
	350	700	350	700		350	700		350	700		
$d = 0.30$	0.9984	0.9992	1.0488	1.0491		0.9982	0.9990		0.9485	0.9489		
$d = 0.60$	0.9987	0.9997	1.0491	1.0497		0.9984	0.9995		0.9488	0.9495		
$d = 1.00$	0.9990	1.0002	1.0494	1.0502		0.9987	1.0000		0.9490	0.9500		

Table 15: Mean estimate: Small breaks. This table reports the mean of the full-sample LCM estimates for different predictive models. Specifically, the estimator is implemented using trimming $\ell = \lfloor n^{\nu} \rfloor$ and bandwidth $m = \lfloor n^{\kappa} \rfloor$, for $(\nu, \kappa) = (0.2, 0.7)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_0 describes the null hypothesis without parameter instability. H_4 , H_5 , H_6 describes three predictive models with small breaks. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

Standard deviation of LCM Estimate: Small breaks												
	No Cointegration: $b=0$											
	$\mathcal{H}_0, n =$			\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$		\mathcal{H}_6 , $n=$				
	350	700	350	700	350	700	350	700				
$d = 0.30$	0.0471	0.0366	0.0483	0.0376	0.0486	0.0375	0.0482	0.0372				
$d = 0.60$	0.0471	0.0366	0.0483	0.0376	0.0486	0.0375	0.0482	0.0372				
$d = 1.00$	0.0471	0.0366	0.0483	0.0376	0.0486	0.0375	0.0482	0.0372				
	Cointegration: $b = d$											
	\mathcal{H}_0 , $n=$			\mathcal{H}_4 , $n=$		\mathcal{H}_5 , $n=$		\mathcal{H}_6 , $n=$				
	350	700	350	700	350	700	350	700				
$d = 0.30$	0.0388	0.0295	0.0403	0.0308	0.0408	0.0307	0.0401	0.0302				
$d = 0.60$	0.0335	0.0241	0.0352	0.0257	0.0359	0.0257	0.0350	0.0252				
$d = 1.00$	0.0291	0.0185	0.0309	0.0206	0.0319	0.0210	0.0308	0.0202				

Table 16: Standard deviation: Small breaks. This table reports the standard deviation of the full-sample LCM estimates for different predictive models. Specifically, the estimator is implemented using trimming $\ell = |n^{\nu}|$ and bandwidth $m = \lfloor n^{\kappa} \rfloor$, for $(\nu, \kappa) = (0.2, 0.7)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . H₀ describes the null hypothesis without parameter instability. \mathcal{H}_4 , \mathcal{H}_5 , \mathcal{H}_6 describes three predictive models with small breaks. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

Bias of LCM-based Break Fraction Estimate: Small Breaks												
	No Cointegration: $b = 0$											
	\mathcal{H}_4 , $n=$			\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$					
	350	700		350	700		350	700				
$d = 0.30$	0.1026	0.0712		-0.0023	0.0055		-0.1062	-0.0591				
$d = 0.60$	0.1026	0.0712		-0.0023	0.0055		-0.1062	-0.0591				
$d = 1.00$	0.1026	0.0712		-0.0023	0.0055		-0.1062	-0.0591				
	Cointegration: $b = d$											
	\mathcal{H}_4 , $n=$			\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$					
	350	700		350	700		350	700				
$d = 0.30$	0.0772	0.0363		-0.0270	-0.0167		-0.1231	-0.0769				
$d = 0.60$	0.0572	0.0242		-0.0324	-0.0228		-0.1218	-0.0733				
$d = 1.00$	0.0500	0.0136		-0.0342	-0.0175		-0.1300	-0.0564				

Table 17: Break fraction bias: Small breaks. This table reports the bias of the break fraction estimates from the LCM-based sup-Wald test in (19). Specifically, the test is implemented using trimming $\ell = |n^{\nu}|$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . \mathcal{H}_4 , \mathcal{H}_5 , \mathcal{H}_6 describes three predictive models with small breaks, whose break fractions are 0.25, 0.50 and 0.75, respectively. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

RMSE of LCM-based Break Fraction Estimate: Small Breaks										
	No Cointegration: $b = 0$									
	\mathcal{H}_4 , $n=$			\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$			
	350	700		350	700	350	700			
$d = 0.30$	0.1992	0.1643		0.1532	0.1286	0.2061	0.1449			
$d = 0.60$	0.1992	0.1643		0.1532	0.1286	0.2061	0.1449			
$d = 1.00$	0.1992	0.1643		0.1532	0.1286	0.2061	0.1449			
	Cointegration: $b = d$									
	\mathcal{H}_4 , $n=$			\mathcal{H}_5 , $n=$			\mathcal{H}_6 , $n=$			
	350	700		350	700	350	700			
$d = 0.30$	0.1683	0.1051		0.1396	0.1056	0.2201	0.1641			
$d = 0.60$	0.1373	0.0864		0.1307	0.0926	0.2146	0.1587			
$d = 1.00$	0.1273	0.0604		0.1282	0.0783	0.2220	0.1373			

Table 18: Break fraction RMSE: Small breaks. This table reports the bias of the break fraction estimates from the LCM-based sup-Wald test in (19). Specifically, the test is implemented using trimming $\ell = |n^{\nu}|$, bandwidth $m = \lfloor n^{\kappa} \rfloor$ and a corresponding covariance estimator bandwidth $m_G = \lfloor n^{\kappa_G} \rfloor$, for $(\nu, \kappa, \kappa_G) = (0.2, 0.7, 0.9)'$. As described in the main text, $\{y_t, x_{t-1}\} \in I(d)$, with x_{t-1} predicting y_t . H₄, H₅, H₆ describes three different predictive models with small breaks, whose break fractions are 0.25, 0.50 and 0.75, respectively. The predictive model is specified with $(b = d)$ and without $(b = 0)$ cointegration. The sample size is either $n = 350$ or $n = 700$, and the nominal test size is 5%. Simulations are implemented with 1000 replications.

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