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# 1 Introduction

When modelling prices of financial assets, modelling the volatility process itself plays a crucial role. In the derivatives world, log-prices are often modelled as continuous semi-martingales. For a given asset with log-price  $Y_t$ , such a process takes the form

$$dY_t = \mu_t dt + \sigma_t dW_t \quad (1)$$

where  $\mu_t$  is a drift term and  $W_t$  is a one-dimensional Brownian motion. The term  $\sigma_t$  represents the volatility process, and it is a very important ingredient in the model [Gatheral et al. \[2018\]](#). However, modelling the volatility process itself is also a complex task. Different approaches to estimate  $\sigma_t$  is used in the literature. In the simplest models, the volatility is either constant or a deterministic function of time. In more popular stochastic volatility models, the volatility process is modelled as a stochastic process itself. In notable models such as the Hull and White model, the Heston model, and the SABR model, the volatility  $\sigma_t$  is modelled as a continuous Brownian semi-martingale [Gatheral et al. \[2018\]](#). The stochastic volatility models seem way more realistic but generated prices from these models are in many cases not consistent with observed prices. (33, 46)

Various authors have proposed other kinds of models to model the volatility. A popular one is stochastic volatility models driven by fractional Brownian motion. A well-known example of such a fractional stochastic volatility model is the one proposed by [Comte and Renault \[1998\]](#) who models the dynamics of the volatility  $\sigma_t$  of an asset as:

$$Y_t = \ln \sigma_t, \quad dY_t = -\gamma Y_t dt + \theta dB_t^H \quad (2)$$

where  $B^H$  is a fractional Brownian motion with Hurst exponent  $H$ . The model was first introduced to model long range dependence effect observed in financial time series. The long range dependence is modelled by choosing  $1 > H > \frac{1}{2}$  [Comte and Renault \[1998\]](#). In recent literature starting with [Gatheral et al. \[2018\]](#), it has been suggested to use the fractional stochastic volatility models with  $H < \frac{1}{2}$  for modelling volatility. Processes driven by a fractional Brownian motion with  $H < \frac{1}{2}$  are referred to as 'rough processes' since these fractional Brownian motion have trajectories rougher than a standard Brownian motion [Cont and Das \[2024\]](#), and [Gatheral et al. \[2018\]](#) concludes that volatility is rough. It is important to note that [Gatheral et al. \[2018\]](#) unlike previous literature rely on the behaviour of volatility estimators over short intraday time scales in order to assess the 'roughness'.

[Cont and Das \[2024\]](#) challenges the conclusions of [Gatheral et al. \[2018\]](#) and similar articles concluding that volatility is rough. [Cont and Das \[2024\]](#) simulates from models where the true spot volatility is known and shows that measures of roughness for realized volatility based on the data from these simulations are in many cases much rougher than those of the underlying true spot

volatility. This difference solely lies in the estimation error, and challenges the use of high-frequency volatility estimators when measuring the roughness.

In this thesis, we will take a similar approach as in [Cont and Das \[2024\]](#) and investigate further if volatility appears to be rough even when the true model does not exhibit rough behaviour.

## 2 Fractional Brownian motion

A fractional Brownian motion (fBm)  $(B_t^H)_{t \in \mathbb{R}}$  with Hurst parameter  $H \in (0, 1)$  is a centered continuous Gaussian process with covariance function

$$\mathbb{E}[B_t^H B_s^H] = \frac{1}{2} (t^{2H} + s^{2H} - |t - s|^{2H}) \quad (3)$$

for  $s, t \geq 0$ . Note that  $B_t^H$  reduces to an ordinary Brownian motion for  $H = \frac{1}{2}$  [Dieker and Mandjes \[2003\]](#). The incremental process of a fractional Brownian motion is called fractional Gaussian noise, and it is a stationary discrete-time process. We define the fractional Gaussian noise  $X = \{X_k : k = 0, 1, \dots\}$  by

$$X_k := B_{t_{k+1}}^H - B_{t_k}^H. \quad (4)$$

With this definition, we immediately see that every  $X_k$  is normally distribution with mean 0 and variance

$$\begin{aligned} \mathbb{E}[(B_{t_{k+1}}^H - B_{t_k}^H)^2] &= \text{Var}(B_{t_{k+1}}^H) + \text{Var}(B_{t_k}^H) - 2\text{Cov}(B_{t_{k+1}}^H, B_{t_k}^H) \\ &= t_{k+1}^{2H} + t_k^{2H} - (t_{k+1}^{2H} + t_k^{2H} - (t_{k+1} - t_k)^{2H}) = (t_{k+1} - t_k)^{2H} \end{aligned}$$

using the covariance function (3). Thus, the fractional Gaussian noise is standard normal distributed when the time step is 1.

### 2.1 Simulation of fractional Brownian motion

Throughout this thesis we will be simulating from the fractional Brownian motion by using a spectral method which can be used for stationary processes. The idea is to analyse the stochastic process in the spectral domain rather than the time domain. The spectral density is computed as follows for frequencies  $-\pi \leq \lambda \leq \pi$ :

$$f(\lambda) := \sum_{j=-\infty}^{\infty} \gamma(j) \exp(ij\lambda) \quad (5)$$

where  $\gamma(\cdot)$  represent the autocovariance function. It can be shown that the spectral density of fractional Gaussian noise is given by

$$f(\lambda) = 2 \sin(\pi H) \Gamma(2H + 1) (1 - \cos \lambda) [|\lambda|^{-2H-1} + B(\lambda, H)] \quad (6)$$

where  $\Gamma(\cdot)$  denotes the Gamma function and

$$B(\lambda, H) := \sum_{j=1}^{\infty} ((2\pi j + \lambda)^{-2H-1} + (2\pi j - \lambda)^{-2H-1}) \quad (7)$$

for  $-\pi \leq \lambda \leq \pi$ . The infinite sum makes direct numerical evaluation almost impossible. However, a useful approximation of the sum is made by [Paxson \[1997\]](#). They show that by using

$$\tilde{B}_3(\lambda, H) := \sum_{j=1}^3 \left( (a_j^+)^{-2H-1} + (a_j^-)^{-2H-1} \right) + \frac{(a_3^+)^{-2H} + (a_3^-)^{-2H} + (a_4^+)^{-2H} + (a_4^-)^{-2H}}{8H\pi}$$

where  $a_j^\pm = 2\pi j \pm \lambda$ ,  $f(\lambda)$  is approximated quite well.

Now, consider a stationary Gaussian discrete-time process  $X = \{X_n : n = 0, \dots, N-1\}$  where  $N$  is the required sample size. The spectral theorem states that it can be represented in terms of the spectral density as

$$X_n \stackrel{d}{=} \int_0^\pi \sqrt{\frac{f(\lambda)}{\pi}} \cos(n\lambda) dB_1(\lambda) - \int_0^\pi \sqrt{\frac{f(\lambda)}{\pi}} \sin(n\lambda) dB_2(\lambda) \quad (8)$$

where  $B_1$  and  $B_2$  are two mutually independent Brownian motions [Dieker and Mandjes \[2003\]](#). We wish to approximate (8). The integrand is replaced by a simpler function. Fix some integer  $l$  and set  $t_k = \pi k/l$  for  $k = 0, \dots, l$ . Now, define a simple function  $\xi_n^{(l)}$  on  $[0, \pi]$  for  $0 \leq n \leq N-1$  by

$$\xi_n^{(\ell)}(\lambda) = \sqrt{\frac{f(t_1)}{\pi}} \cos(nt_1) \mathbf{1}_{\{0\}}(\lambda) + \sum_{k=0}^{l-1} \sqrt{\frac{f(t_{k+1})}{\pi}} \cos(nt_{k+1}) \mathbf{1}_{(t_k, t_{k+1}]}(\lambda). \quad (9)$$

The first integral in (8) can be approximated by  $\int_0^\pi \xi_n^{(\ell)}(\lambda) dB_1(\lambda)$ . Since  $\xi_n^{(\ell)}$  is a simple function the stochastic integral can be computed as

$$\int_0^\pi \xi_n^{(l)}(\lambda) dB_1(\lambda) = \sum_{j=0}^{\ell-1} \xi_n^{(l)}(B(t_{j+1}) - B(t_j)). \quad (10)$$

Thus, we obtain

$$\begin{aligned} \int_0^\pi \xi_n^{(l)}(\lambda) dB_1(\lambda) &= \sum_{j=0}^{\ell-1} \sum_{k=0}^{\ell-1} \sqrt{\frac{f(t_{k+1})}{\pi}} \cos(nt_{k+1}) \mathbf{1}_{(t_k, t_{k+1}]}(B_1(t_{j+1}) - B_1(t_j)) \\ &= \sum_{k=0}^{l-1} \sqrt{\frac{f(t_{k+1})}{\pi}} \cos(nt_{k+1}) (B_1(t_{k+1}) - B_1(t_k)) \\ &= \sum_{k=0}^{\ell-1} \sqrt{\frac{f(t_{k+1})}{\pi}} \cos(nt_{k+1}) U_k^{(0)} \sqrt{\frac{\pi}{l}} = \sum_{k=0}^{\ell-1} \sqrt{\frac{f(t_{k+1})}{\ell}} \cos(nt_{k+1}) U_k^{(0)} \end{aligned}$$

where  $U_k^{(0)}$  is an i.i.d. standard normal random variable for  $k = 0, \dots, \ell - 1$ . The  $U_k^{(0)} \sqrt{\pi/\ell}$  represent the Brownian motion increments which per definition are normally distributed with mean 0 and variance  $t_{k+1} - t_k$ .

The second integral in (8) can be approximated in a similar way by replacing the cosine terms with sine terms. Thus, we obtain the following approximation of  $X_n$ :

$$\hat{X}_n^{(\ell)} := \sum_{k=0}^{\ell-1} \sqrt{\frac{f(t_{k+1})}{\ell}} \left( \cos(nt_{k+1}) U_k^{(0)} - \sin(nt_{k+1}) U_k^{(1)} \right). \quad (11)$$

The two vectors  $U^{(0)}$  and  $U^{(1)}$  are mutually independent since  $B_1$  and  $B_2$  are independent as well. In order to calculate  $\hat{X}_n^{(\ell)}$  efficiently we will be using the fast Fourier transform (FFT). To this end, we define the sequence  $(a_k)_{k=0, \dots, 2\ell-1}$  by

$$a_k := \begin{cases} 0 & k = 0; \\ \frac{1}{2} \left( U_{k-1}^{(0)} + i U_{k-1}^{(1)} \right) \sqrt{\frac{f(t_k)}{\ell}} & k = 1, \dots, \ell - 1; \\ U_{k-1}^{(0)} \sqrt{\frac{f(t_k)}{\ell}} & k = \ell; \\ \frac{1}{2} \left( U_{2\ell-k-1}^{(0)} - i U_{2\ell-k-1}^{(1)} \right) \sqrt{\frac{f(t_{2\ell-k})}{\ell}} & k = \ell + 1, \dots, 2\ell - 1. \end{cases}$$

It is shown in appendix that the Fourier transform of  $a_k$  is indeed real and equals  $\hat{X}_n^{(\ell)}$ . From this approximated fractional Gaussian noise we can generate the fBm.

[Dieker and Mandjes \[2003\]](#) shows that the finite-dimensional distributions of  $\hat{X}^{(\ell)}$  converge in probability to the corresponding finite-dimensional distributions of  $X$  as  $\ell \rightarrow \infty$ . The rate of convergence is, however, quite slow. Therefore, we will be using a  $\ell \geq 30000$  when simulating from the fBm by the spectral method in this thesis in order to make sure that our simulations are reliable. Figure 1 shows three simulations of fractional Brownian motions with different Hurst parameters generated by the spectral method.

### 3 Estimating the roughness of volatility

Determining the roughness of volatility plays a crucial role in order to make proper model specification and design estimators. In practice when estimating the roughness, we observe only a single price path, and we need to measure the roughness of this path. Estimating the roughness from discrete observation is not an easy task and multiple ways of estimating the roughness exist in the literature. In this thesis, we will be working with high frequency data, and we will introduce three different roughness estimators that works on high frequency

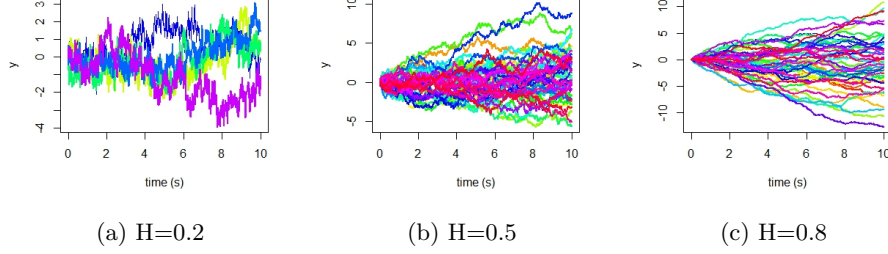


Figure 1: Fractional Brownian motions generated by the spectral method.

data. In section 4 we will conduct numerical experiments, and we will compare the three estimators.

### 3.1 Roughness index estimator via $p$ -th variation

In this section, we will describe the roughness index estimator via normalized  $p$ -th variation which was introduced in [Cont and Das \[2024\]](#). We will closely follow the theory and methodology behind the estimator described in [Cont and Das \[2024\]](#).

We will be working with partitions of our time interval. Consider a sequence of partitions  $\pi = (\pi^n)_{n \geq 1}$  of  $[0, T]$  where

$$\pi^n = \left(0 = t_0^n < t_1^n < \dots < t_{N(\pi^n)}^n = T\right)$$

represents observation times at frequency  $n$ . Here  $N(\pi^n)$  denotes the number of intervals in the partition  $\pi^n$ . For the time partitions we will be working with in this thesis we will assume that

$$|\pi^n| := \sup_{i=1, \dots, N(\pi^n)} |t_i^n - t_{i-1}^n| \xrightarrow{n \rightarrow \infty} 0. \quad (12)$$

That is, the size of the largest interval of  $\pi^n$  will converge to 0 as  $n \rightarrow \infty$ .

We will now define the concept of  $p$ -th variation along a sequence of partitions  $(\pi^n)_{n \geq 1}$ . The definition follows the definition from [Cont and Perkowski \[2019\]](#).

**Definition 1.** ( *$p$ -th variation along a sequence of partitions*)  $x \in C^0([0, T], \mathbb{R})$  has finite  $p$ -th variation along the sequence of partitions  $\pi = (\pi^n, n \geq 1)$  if there exists a continuous increasing function  $[x]_\pi^{(p)} : [0, T] \rightarrow \mathbb{R}_+$  such that

$$\forall t \in [0, T], \quad \sum_{\{t_j^n, t_{j+1}^n\} \in \pi^n : t_j^n \leq t} |x(t_{j+1}^n) - x(t_j^n)|^p \xrightarrow{n \rightarrow \infty} [x]_\pi^{(p)}(t). \quad (3)$$

If this property holds, then the convergence in (3) is uniform. We call  $[x]_\pi^{(p)}$  the  $p$ -th variation of  $x$  along the sequence of partitions  $\pi$ . We denote  $V_\pi^p([0, T], \mathbb{R})$  the set of all continuous paths with finite  $p$ -th variation along  $\pi$ .

We further define the concepts of variation index and roughness index of a path following Cont and Das [2024].

**Definition 2** (Variation index). *The variation index of a path  $x$  along a partition sequence  $\pi$  is defined as the smallest  $p \geq 1$  for which  $x$  has finite  $p$ -th variation along  $\pi$ :*

$$p^\pi(x) = \inf \{p \geq 1 : x \in V_\pi^p([0, T], \mathbb{R})\}.$$

**Definition 3** (Roughness index). *The roughness index of a path  $x$  (along  $\pi$ ) is defined as*

$$H^\pi(x) = \frac{1}{p^\pi(x)}.$$

When the underlying sequence of partitions is clear, we will simply denote these indexes as  $p(x)$  and  $H(x)$ .

For a real-valued stochastic process  $X : [0, T] \times \Omega \rightarrow \mathbb{R}$  the variation index (hence also the roughness index)  $p^\pi(X(., \omega))$  of each sample path  $X(., \omega)$  may in principle be different. However, there are many important classes of stochastic processes which have an almost-sure roughness index. This holds for example for a fractional Brownian motion where the roughness index matches with the corresponding Hurst parameter Cont and Das [2024].

When estimating roughness from empirical data based on discrete observations, using the  $p$ -th variation directly is difficult since it involves checking convergence to an unknown limit. To this end, we introduce the concept of normalized  $p$ -th variation.

**Definition 4** (Normalized  $p$ -th variation along a sequence of partitions). *Let  $\pi$  be a sequence of partitions of  $[0, T]$  with mesh  $|\pi^n| \rightarrow 0$  and  $\pi^n = \{0 = t_1^n < t_2^n < \dots < t_{N(\pi^n)}^n = T\}$ .  $x \in V_\pi^p([0, T], \mathbb{R})$  is said to have normalized  $p$ -th variation along  $\pi$  if there exists a continuous function  $w(x, p, \pi) : [0, T] \rightarrow \mathbb{R}$  such that:*

$$\forall t \in [0, T], \quad \sum_{\pi^n \cap [0, t]} |x(t_{j+1}^n) - x(t_j^n)|^p \rightarrow w(x, p, \pi)(t).$$

(Cont and Das 2022) shows that for a large class of functions with  $p$ -th variation, the normalized  $p$ -th variation exists and is linear. Furthermore, Cont and Das [2024] shows that the normalized  $p$ -th variation is a 'sharp' statistic meaning that if a function has finite  $p$ -th variation then for all  $q \neq p$  the normalized  $p$ -th variation is either zero or infinite. Additionally, it is shown that a fractional

Brownian motion has normalized  $p$ -th variation along the dyadic partition  $\mathbb{T} = (\mathbb{T}^n)_{n \geq 1}$  almost-surely.

### 3.1.1 Estimating the roughness from discrete observations

With these concepts defined we now proceed to explaining how roughness can be estimated from discrete observations. Given discrete observations on a refining partition  $\pi^L$  we define the 'normalized  $p$ -th variation statistic' which is the discrete counterpart of the normalized  $p$ -th variation:

$$W(L, K, \pi, p, t, X) := \sum_{\pi^K \cap [0, t]} \frac{|X(t_{i+1}^K) - X(t_i^K)|^p}{\sum_{\pi^L \cap [t_i^K, t_{i+1}^K]} |X(t_{j+1}^L) - X(t_j^L)|^p} \times (t_{i+1}^K - t_i^K). \quad (13)$$

This definition involves two frequencies  $K$  and  $L \gg K$ . We can consider  $L$  as the sample frequency while  $K$  is a block frequency such that  $\pi^K$  is a subpartition of  $\pi^L$ . Thus, we are grouping the sample size  $L$  into  $K$  many groups where each group contains exactly  $\frac{L}{K}$  consecutive points. It can be seen that the normalized  $p$ -th variation statistic converges to the normalized  $p$ -th variation as  $L$  and  $K$  increase. That is,

$$\lim_{K \rightarrow \infty} \lim_{L \rightarrow \infty} W(L, K, \pi, p, t, X) = w(x, p, \pi)(t). \quad (14)$$

The variation index estimator  $\hat{p}_{L,K}^\pi(X)$  associated with data sampled on  $\pi^L$  can then be obtained by computing  $W(L, K, \pi, p, t, X)$  for different values of  $p$  such that the following equation can be solved for  $p_{L,K}^\pi(X)$

$$W(L, K, \pi, \hat{p}_{L,K}^\pi(X), t, X) = T. \quad (15)$$

The corresponding roughness index estimator is then defined as

$$\hat{H}_{L,K}^\pi(X) = \frac{1}{\hat{p}_{L,K}^\pi(X)}.$$

If the underlying dataset and partition is clear we will also denote these estimators as  $\hat{p}_{L,K}$  and  $\hat{H}_{L,K}$ . Asymptotic properties of these estimators under high-frequency sampling are studied in (cont and das 2022).

### 3.1.2 Behaviour of the roughness estimator based on simulations

We will now study the behaviour of the roughness estimator based on high-frequency simulated data. We will be simulating from a fractional Brownian motion and compute the roughness estimator based on the data. In this simulation section we will be using a uniform partition of the time interval  $[0, 1]$  with

$$\pi^n = \left( 0 < \frac{1}{n} < \frac{2}{n} < \dots < 1 \right).$$



We will be simulating from four fractional Brownian motions with four different Hurst exponents  $H = \{0.1, 0.3, 0.5, 0.8\}$  in order to investigate the accuracy of the roughness estimator  $\hat{H}_{L,K}^\pi(X)$ . First, we plot  $\log(W(L = 300 \times 300, K = 300, \pi, p, t = 1, X = B^H))$  against  $H = 1/p$  to visualize the estimation. The results are presented in Figure 2. The solid black line is the value of  $\log(W(L = 300 \times 300, K = 300, \pi, p, t = 1, X = B^H))$  for different values of  $1/p$ . The blue vertical line represents the estimated roughness index whereas the black dotted vertical line represent the true Hurst parameter. The blue horizontal line represents  $W(L, K, \pi, p, t, X) = T$ . The estimated roughness index  $\hat{H}_{L,K}$  is the crossing between  $\log(W(L = 300 \times 300, K = 300, \pi, p, t = 1, X = B^H))$  and  $W(L, K, \pi, p, t, X) = T$ .

In Figure 3, we have created histograms of the estimator  $\hat{H}_{L,K}$  again by

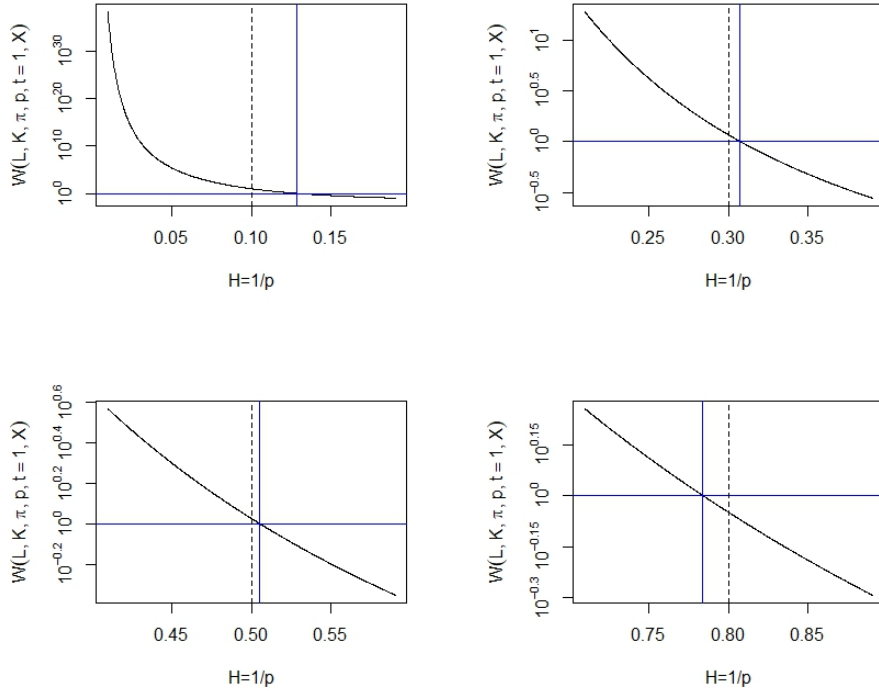


Figure 2: Log-scale plot of the normalized  $p$ -th variation statistic for fBm with Hurst parameter  $H = \{0.1, 0.3, 0.5, 0.8\}$ . The black solid line represents the value of  $\log(W(L = 300 \times 300, K = 300, \pi, p, t = 1, X = B^H))$  plotted against  $H = 1/p$ . The blue vertical line represents  $\hat{H}_{L,K}$  using the normalized  $p$ -th variation statistic with  $L = 300 \times 300$  and  $K = 300$ . The vertical black dotted line represents the true Hurst parameter.

using  $W(L = 300 \times 300, K = 300, \pi, p, t = 1, X = B^H)$  from 150 independent simulations. In addition to this, Table 1 provides the summary statistics for the estimated roughness index  $\hat{H}_{L,K}$  shown in the histograms. Based on Figure 2 and 3, and Table 1, we conclude that the estimated roughness index  $\hat{H}_{L,K}$  seems to be fairly accurate when used on a data set with length  $L = 300 \times 300$  generated from a fractional Brownian motion. All the estimated  $\hat{H}_{L,K}$  are within a distance 0.05 from the true Hurst parameter, and the mean and median of the estimated roughness are in all cases very close to the true Hurst parameter indicating that our estimator is quite accurate. Only exception is for the fBm with Hurst parameter  $H = 0.8$  where even the upper quartile of  $\hat{H}_{L,K}$  is smaller than the true Hurst parameter as seen in Table 1. This suggest that the estimated roughness index might be biased downwards in the case of  $H = 0.8$ . The estimate is, however, still quite accurate even for  $H = 0.8$ .

H	Min.	Lower quartile	Median	Mean	Upper quartile	Max.
0.1	0.0622	0.0944	0.1021	0.1031	0.1149	0.1402
0.3	0.2746	0.2954	0.3012	0.3005	0.3052	0.3177
0.5	0.4743	0.4954	0.5004	0.4997	0.5041	0.5239
0.8	0.7657	0.7797	0.7856	0.7865	0.7937	0.8239

Table 1: Summary of statistics for estimated roughness index  $\hat{H}_{L,K}$  from 150 independent simulations of fBm with  $L = 300 \times 300$  and  $K = 300$ .

In Figure 4 we have used  $L = 2000 \times 2000$  and  $K = 2000$  to make similar plots to those in Figure 2 and Figure 3 for a fractional Brownian motion with Hurst parameter  $H = 0.1$ . Table 2 provides the summary statistics for the estimated roughness index  $\hat{H}_{L,K}$  corresponding to the histogram in Figure 4. We observe that the results are similar to those we obtained for  $L = 300 \times 300$  and  $K = 300$  but that the estimated roughness index  $\hat{H}_{L,K}$  seems to be even more accurate as seen in Figure 4 and Table 2. This is not surprising since we know that the normalized  $p$ -th variation statistic converges to the normalized  $p$ -th variation as  $L$  and  $K$  increase. However, increasing  $L$  and  $K$  is also computationally more demanding. The accuracy of  $\hat{H}_{L,K}$  when using  $L = 300 \times 300$  and  $K = 300$  is satisfactory for many of our purposes. Therefore, we will be using at least  $L = 300 \times 300$  and  $K = 300$  for the remainder of the thesis.

H	Min.	Lower quartile	Median	Mean	Upper quartile	Max.
0.1	0.0892	0.0963	0.0996	0.0998	0.1031	0.1145

Table 2: Summary of statistics for estimated roughness index  $\hat{H}_{L,K}$  from 150 independent simulations of fBm with  $H = 0.1$ ,  $L = 2000 \times 2000$  and  $K = 2000$ .

We further investigate how the choice of  $K \ll L$  affects the estimated roughness index  $\hat{H}_{L,K}$ . In Figure 5, we have plotted the estimated roughness  $\hat{H}_{300 \times 300, K}$

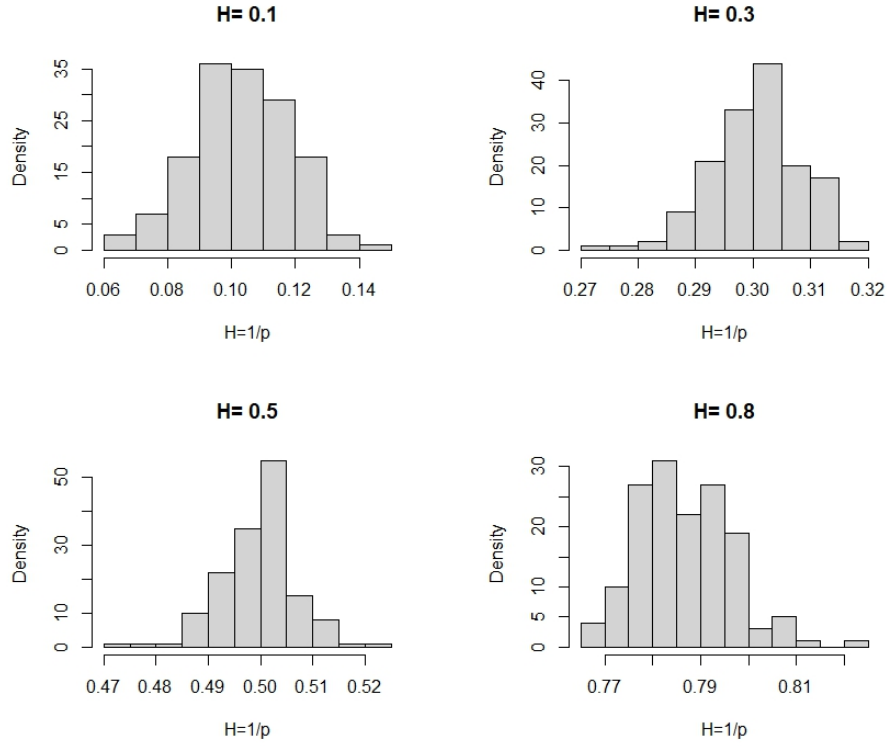


Figure 3: Histogram of estimated roughness index  $\hat{H}_{L,K}$  with  $L = 300 \times 300$  and  $K = 300$  generated from 150 independent simulation of fractional Brownian motions with Hurst parameter  $H = \{0.1, 0.3, 0.5, 0.8\}$ .

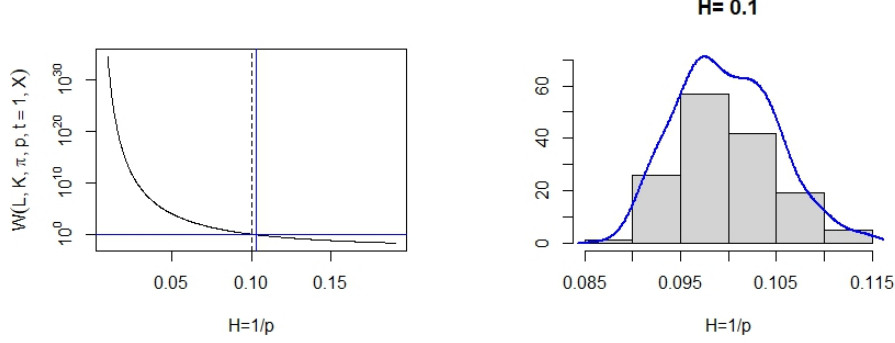


Figure 4: Simulation results for fractional Brownian motion with Hurst parameter  $H = 0.1$ . **Left:** The value of  $\log(W(L = 2000 \times 2000, K = 2000, \pi, p, t = 1, X = B^H))$  plotted against  $H = 1/p$  in black. The blue vertical line represents the estimated roughness  $\hat{H}_{L,K}$  using  $L = 2000 \times 2000$  and  $K = 2000$ . The vertical black dotted line represents the true Hurst parameter. **Right:** Histogram of estimated roughness index  $\hat{H}_{L,K}$  using  $L = 2000 \times 2000$  and  $K = 2000$  generated by 150 independent simulations of fBm with Hurst parameter  $H = 0.1$ . The blue line represents a kernel estimator for density.

for a fractional Brownian motion with Hurst parameter  $H = 0.1$  for different values of  $K$  with a fixed  $L = 300 \times 300$ . Note that when  $\frac{L}{K}$  is not an integer, the  $K$  many groups from the definition of normalized  $p$ -th variation statistic (13) do not contain exactly  $\frac{L}{K}$  consecutive points. The code is implemented such that each group will contain either  $\lceil \frac{L}{K} \rceil$  or  $\lfloor \frac{L}{K} \rfloor$  consecutive points. Figure 5 shows that when  $K$  is too low, the estimator  $\hat{H}_{300 \times 300, K}$  seems to be underestimating the true Hurst parameter whereas when  $K$  is too high the Hurst parameter is overestimated. For  $K \approx \sqrt{L}$  the estimated roughness index is quite consistent and close to the true Hurst parameter  $H = 0.1$ . It is thus natural to use  $K = \sqrt{L}$  since  $\frac{L}{K}$  will be an integer and the estimator  $\hat{H}_{L,K}$  seems to be accurate and consistent in that range.

### 3.2 Instantaneous volatility and realized volatility

As mentioned in (1) log-prices are often modelled as continuous semi-martingales where the derivative of the log-price  $Y_t$  takes the form

$$dY_t = \mu_t dt + \sigma_t dW_t.$$

The term  $\sigma_t$  represents the volatility process, and  $\sigma_t$  is also called instantaneous volatility or spot volatility. In stochastic volatility models,  $\sigma_t$  is represented as a stochastic process itself often driven by a fractional process. Contrary to prices

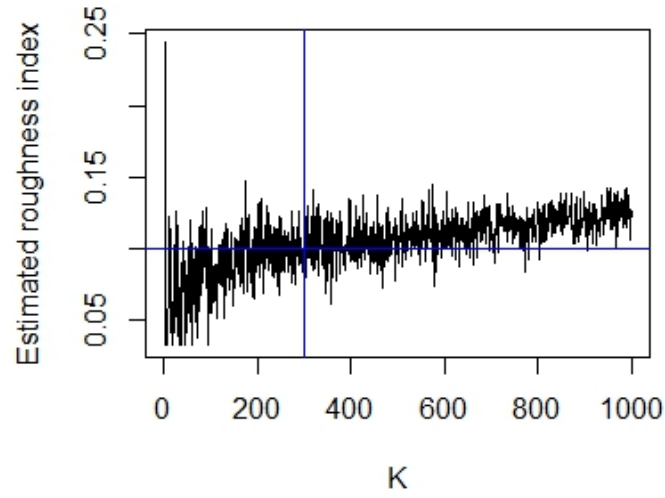


Figure 5: The solid black line represents the estimated roughness index  $\hat{H}_{300 \times 300, K}$  plotted against different values of  $K$  for a simulation from a fBm with Hurst parameter  $H = 0.1$ . The blue vertical line represents  $K = \sqrt{L} = 300$  whereas the blue horizontal line represents the true Hurst parameter  $H = 0.1$ .

of an asset, instantaneous volatility cannot be directly observed and needs to be estimated from prices.

In a practical situation, the price of the asset  $S_t$  at time  $t$  is usually observed over a non-uniform time grid of  $[0, T]$ :

$$\pi^n = \left(0 = t_0^n < t_1^n < \dots < t_{N(\pi^n)}^n = T\right).$$

For high-frequency data the assumption (12) is often assumed. That is,  $|\pi^n| \rightarrow 0$  as  $n$  increases where  $n$  can be thought of as a sampling frequency i.e. the number of samples per second (or per other unit).

For  $Y_t = \log(S_t)$  the instantaneous volatility  $\sigma_t$  can be recovered by

$$\sigma_t^2 = \frac{d}{dt} \lim_{n \rightarrow \infty} \sum_{\pi^n \cap [0, t]} (Y(t_{i+1}^n) - Y(t_i^n))^2 = \lim_{n \rightarrow \infty} RV_t^2(\pi^n)^2$$

where the realized variance along the sampling grid  $\pi^n$  is defined as

$$RV_t(\pi^n)^2 = \sum_{\pi^n \cap [0, t]} (Y(t_{i+1}^n) - Y(t_i^n))^2.$$

The realized volatility over time interval  $[t, t + \Delta]$  along the sampling grid  $\pi^n$  is then defined as

$$RV_{t, t+\Delta}(\pi^n) = \sqrt{\sum_{\pi^n \cap [t, t+\Delta]} (Y(t_{i+1}^n) - Y(t_i^n))^2}. \quad (16)$$

If the price  $S_t$  follows a stochastic volatility model as, it is known that the realized variance converges to the quadratic variation of  $Y$  as sampling frequency increases (Jacod and Protter)

$$RV_{t, t+\Delta}(\pi^n) \xrightarrow[n \rightarrow \infty]{\mathbb{P}} \sqrt{\int_t^{t+\Delta} \sigma_u^2 du}.$$

Hence, along a single price path observed at high-frequency, we can compute the realized volatility (16) and use this as an indicator of volatility

$$RV_{t, t+\Delta}(\pi^n) \simeq \sqrt{\Delta} \sigma_t.$$

As  $n$  increases the estimation becomes more accurate [Cont and Das \[2024\]](#).

### 3.3 Smoothness of a path by logarithmic regression

In this section we will introduce the smoothness estimator via logarithmic regression used in [Gatheral et al. \[2018\]](#). Consider a volatility process on a time grid  $[0, T]$ . First we pretend that we have access to discrete observations of the spot volatility process with mesh  $\Delta$  such that the observations are

$\sigma_0, \sigma_\Delta, \dots, \sigma_{k\Delta}, \dots$  for  $k \in \{0, \lfloor T/\Delta \rfloor\}$ . Here  $\Delta$  is a natural number and  $\Delta \geq 1$ . If we set  $N = \lfloor T/\Delta \rfloor$ , then for  $q \geq 0$  we define

$$m(q, \Delta) = \frac{1}{N} \sum_{k=1}^N |\log(\sigma_{k\Delta}) - \log(\sigma_{(k-1)\Delta})|^q.$$

Following [Gatheral et al. \[2018\]](#) the main assumption is that for some  $s_q > 0$  and  $b_q > 0$

$$N^{qs_q} m(q, \Delta) \rightarrow b_q \quad (17)$$

as  $\Delta$  tends to zero. Under additional technical assumptions, this essentially means that the volatility process belongs to a Besov smoothness space  $\mathcal{B}_{q,\infty}^{s_q}$  and does not belong to  $\mathcal{B}_{q,\infty}^{s'_q}$  for  $s'_q > s_q$ . Hence,  $s_q$  can be considered as a smoothness parameter. In particular, if  $\log(\sigma_t)$  is a fractional Brownian motion with Hurts parameter  $H$ , then for any  $q \geq 0$  equation (17) holds in probability with  $s_q = H$ . It can further be shown that the sample paths of the process do indeed belong to  $\mathcal{B}_{q,\infty}^{s_q}$ . Thus,  $s_q$  is a measure of the smoothness (or roughness) of the volatility process [Gatheral et al. \[2018\]](#).

The instantaneous volatility can not be directly observed, and exact computations of  $m(q, \Delta)$  is not possible in practice. In order to make use of  $m(q, \Delta)$  we must therefore approximate the true spot volatility. The estimated volatility can be computed by the realized volatility or some variation of realized volatility. In the following we will be using the notation  $m(q, \Delta)$  with the understanding that we are only approximating the true spot volatility. To estimate the smoothness parameter  $s_q$  for each  $q$ , we can compute  $m(q, \Delta)$  for different values of  $\Delta$  and regress  $\log m(q, \Delta)$  against  $\log \Delta$ . Note that for a given  $\Delta$  several  $m(q, \Delta)$  can be computed depending on the starting point. For example, if  $\Delta = 3$  then the starting point can be  $\hat{\sigma}_0, \hat{\sigma}_\Delta$  or  $\hat{\sigma}_{2\Delta}$ . The final measure of  $m(q, \Delta)$  is computed as the average of these values of  $m(q, \Delta)$  with different starting points.

[Gatheral et al. \[2018\]](#) shows that for a given  $q$ ,  $\log m(q, \Delta)$  values regressed against  $\log \Delta$  essentially lie on a straight line. Assuming stationary increments of the log-volatility, this implies that the increments fulfil the scaling property

$$\mathbb{E} [|\log(\sigma_\Delta) - \log(\sigma_0)|^q] = b_q \Delta^{\zeta_q}$$

where  $\zeta_q = qs_q > 0$  is the slope of the line associated to  $q$ . This result uses that  $m(q, \Delta)$  can be seen as the empirical counterpart of  $\mathbb{E} [|\log(\sigma_\Delta) - \log(\sigma_0)|^q]$ . Furthermore, [Gatheral et al. \[2018\]](#) shows that  $s_q$  seems to not depend on  $q$ , and plotting  $\zeta_q$  against  $q$  shows  $\zeta_q \sim qs_q$ .

Thus, we can compute an estimate of the smoothness (or roughness) of a volatility process by first regressing  $\log m(q, \Delta)$  against  $\log \Delta$  for different values of  $q$ . The slope of the straight line is an estimate of  $\zeta_q$ . Next, we regress  $\zeta_q$  against

$q$ . The slope of the straight line is an estimate of the smoothness parameter  $s_q$ . If  $\log(\sigma_t)$  is a fractional Brownian motion then  $s_q = H$  holds in probability where  $H$  is the Hurst parameter.

In Figure 6 we have replicated the method used in [Gatheral et al. \[2018\]](#) to estimate the roughness index of the volatility of S&P500 using the first 3500 days of the daily realized variance estimates from the Oxford-Man Institute of Quantitative Finance Realized Library for S&P. Our results coincide with the results from [Gatheral et al. \[2018\]](#), and the log-regression yields a straight line for all our values of  $q$ . Our estimated smoothness of the S&P500 volatility is  $H = 0.1421$ .

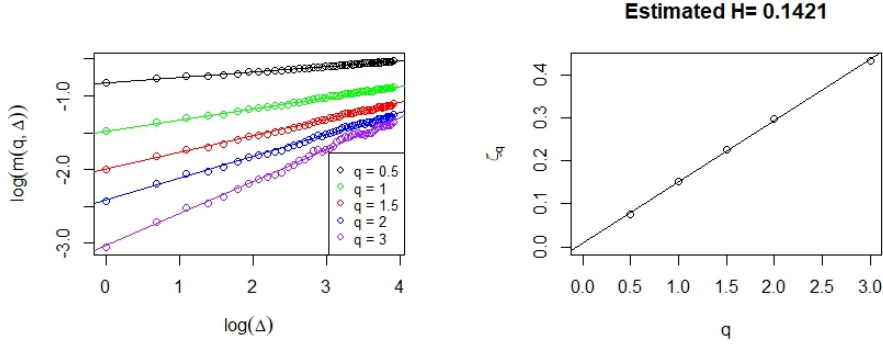


Figure 6: A reproduction of the log-regression method introduced by [Gatheral et al. \[2018\]](#) using the first 3500 days of the daily realized variance estimates from the Oxford-Man Institute of Quantitative Finance Realized Library for S&P. The estimated roughness is  $H = 0.1421$ .

In Figure 7 we have used the roughness estimator from section 3.1 on the same S&P500 volatility data to estimate the roughness. We have used  $W(L = 3500, K = \sqrt{3500}, \pi, p, t = 1, X)$  since the dataset only consist of 3500 days. We have set  $t = 1$  since  $t$  simply works as a scaling parameter. The conclusions are the same for  $t = 1$  as for  $t = 3500$ . The estimated roughness index is  $\hat{H}_{L,K} = 0.1425$ . We immediately observe that the estimated roughness index is very close to the estimated smoothness parameter from the log-regression method  $s_q = 0.1421$ . However, note that  $L = 3500$  is much smaller than  $L = 300 \times 300 = 90000$  for which we have investigated the accuracy of the roughness estimator via normalized  $p$ -th variation statistic. Therefore, the estimated roughness estimator  $\hat{H}_{L,K}$  might not be very precise. The results do however indicate that our two different approaches to estimate the roughness of a volatility process coincides.



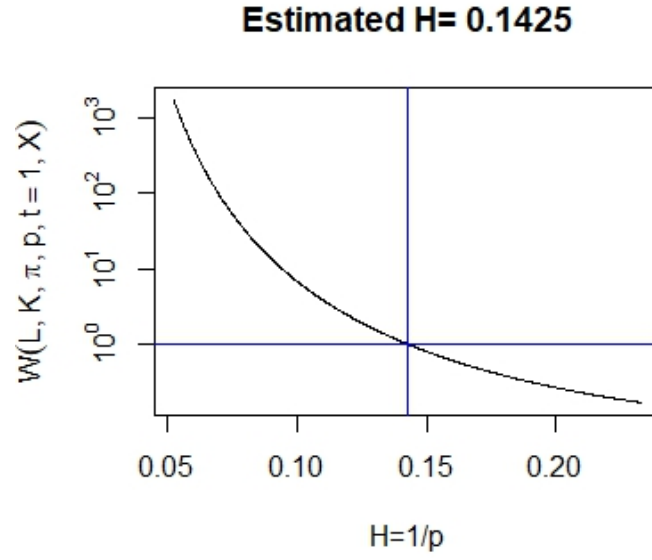


Figure 7: Estimating the roughness of the first 3500 days of S&P500 volatility data from Oxford-Man Institute of Quantitative Finance Realized Library using the roughness estimator from section 3.1 with  $L = 3500$  and  $K = \sqrt{3500}$ . The estimated roughness index is  $\hat{H}_{L,K} = 0.1425$ .

### 3.4 Sequential scale estimator of roughness exponent

The third estimator we will introduce, is the sequential scale estimator of roughness first introduced by [Han and Schied \[2023\]](#). This estimator differs from many other roughness estimators by the fact that it is computed directly from discrete observations of realized variance. That is, the input to the estimator needs to be realized variance and not instantaneous variance or volatility. The two previously described estimators are based on having access to discrete observations of instantaneous volatility  $\sigma_t$ . Realized volatility is then used as an approximation of the true spot volatility. This two step approach can be problematic since the estimation error in the realized volatility might substantially distort the outcome of the final roughness estimation. The sequential scale estimator avoids these possible inaccuracies caused by the estimation errors since it is based directly on realized variance.

[Han and Schied \[2023\]](#) are mainly concerned with stochastic volatility models based on fractional Brownian motions, but many aspects of the approach works in a model-free setting. Let  $x : [0, 1] \leftarrow \mathbb{R}$  be any continuous function. For  $p \geq 1$  the  $p$ -th variation of the function  $x$  along the  $n$ -th dyadic partition is defined as

$$\langle x \rangle_n^{(p)} := \sum_{k=0}^{2^n-1} |x((k+1)2^{-n}) - x(k2^{-n})|^p.$$

The above definition is similar to Definition 1. If there exist a  $R \in [0, 1]$  such that

$$\lim_{n \rightarrow \infty} \langle x \rangle_n^{(p)} = \begin{cases} 0 & \text{for } p > \frac{1}{R} \\ \infty & \text{for } p < \frac{1}{R} \end{cases}$$

we refer to  $R$  as the roughness exponent of  $x$ . The smaller  $R$  the rougher is the path  $x$  and vice verse. Moreover, if  $x$  is a sample path of a fractional Brownian motion, the roughness exponent  $R$  is equal to the Hurst parameter  $H$  [Han and Schied \[2023\]](#).

Now, since only asset prices and their realized variance can be observed in practice, we can consider this as making discrete observations of

$$y(t) = \int_0^t g(x(s)) ds, \quad 0 \leq t \leq 1, \quad (18)$$

where  $g : \mathbb{R} \rightarrow \mathbb{R}$  is sufficiently regular. Here  $y(t)$  correspond to realized variance. For instance, if log-volatility is given by a fractional Ornstein-Uhlenbeck process, we will take  $x$  as a sample path of a fractional Ornstein-Uhlenbeck process and  $g(t) = (e^t)^2 = e^{2t}$ .

For the estimator, it is supposed that for a given  $n \in \mathbb{N}$  we have access to the

discrete observations  $\{y(k2^{-n-2}) : k = 0, \dots, 2^{n+2}\}$  of the realized variance (i.e. the function  $y$  in (18)). Based on these data points, the following coefficients are introduced

$$\vartheta_{n,k} := 2^{3n/2+3} \left( y \left( \frac{4k}{2^{n+2}} \right) - 2y \left( \frac{4k+1}{2^{n+2}} \right) + 2y \left( \frac{4k+3}{2^{n+2}} \right) - y \left( \frac{4k+4}{2^{n+2}} \right) \right), \quad (19)$$

for  $0 \leq k \leq 2^n - 1$ . The estimator for the roughness exponent is then given by

$$\hat{\mathcal{R}}_n(y) := 1 - \frac{1}{n} \log_2 \sqrt{\sum_{k=0}^{2^n-1} \vartheta_{n,k}^2}.$$

Detailed explanation of the rationale behind the estimator can be seen in [Han and Schied \[2023\]](#) where convergence and consistency results are also shown.

The estimator  $\hat{\mathcal{R}}_n(y)$  is not scale-invariant, and [Han and Schied \[2023\]](#) suggest to do scale-invariant modifications of  $\hat{\mathcal{R}}_n(y)$ . Fix  $m \in \mathbb{N}$  with  $n > m$  and fix  $\alpha_0, \dots, \alpha_m \geq 0$  with  $\alpha_0 > 0$ . The sequential scaling factor  $\lambda_n^s$  is then defined as

$$\lambda_n^s := \arg \min_{\lambda > 0} \sum_{k=n-m}^n \alpha_{n-k} \left( \hat{\mathcal{R}}_k(\lambda y) - \hat{\mathcal{R}}_{k-1}(\lambda y) \right)^2.$$

The sequential scaling estimate  $\hat{\mathcal{R}}_n^s(y)$  is then defined as follows

$$\hat{\mathcal{R}}_n^s(y) := \hat{\mathcal{R}}_n(\lambda_n^s y).$$

The sequential scaling factor  $\lambda_n^s$  enforces the convergence of  $\hat{\mathcal{R}}_n(\lambda_n^s y)$ . Thus,  $\hat{\mathcal{R}}_n^s(y)$  will converge faster than  $\hat{\mathcal{R}}_n(y)$ . We will be using the sequential scaling estimate  $\hat{\mathcal{R}}_n^s(y)$  when performing numerical experiments in section 4.

## 4 Numerical Experiments

In this section we will be performing a number of numerical experiments. We will be estimating the roughness index for stochastic models based directly on the instantaneous volatility  $\sigma_t$  and based on the realized volatility by using price trajectories simulated from the models. By doing this we can investigate how estimation errors impact the estimated roughness index. We will be performing these experiments for stochastic models with various degrees of roughness.

### 4.1 Simple stochastic volatility diffusion model

First, we will consider the following stochastic volatility model where the volatility is the modulus of a Brownian motion:

$$dS_t = \sigma_t S_t dB_t \quad \text{with} \quad \sigma_t = |W_t|, \quad (20)$$

where  $S_t$  is the price of the asset and  $B_t$  and  $W_t$  are two independent Brownian motions. For the simulation we use  $S_0 = 1$  and  $T = 1$ . The true roughness of this model is simply  $H = 0.5$ .

We are estimating the realized variance using (16) for 300 consecutive points which corresponds to a 5 minute moving window when the price is updated every second. We estimate the roughness of the volatility process by using the realized volatility and instantaneous volatility as data in our roughness estimators. For every 300 consecutive points we are computing the realized volatility  $RV_{t,t+\delta}(\pi^n)$  as in (16) and rescale the volatility to represents the same time interval as the corresponding instantaneous volatility  $\sigma_t$ . The corresponding instantaneous volatility is then the  $\sigma_t$  for the first of these 300 points. In Figure 8, we have made plots for the realized and instantaneous volatility. The left plot of Figure 8 represents the realized volatility of the price process in black and the instantaneous volatility in red. The middle plot of Figure 8 represents the difference between the realized and instantaneous volatility which is the estimation error. The right plot of Figure 8 represents the log difference between the realized and instantaneous volatility. Figure 8 indicates that the estimation error has a complicated dependence structure. There seems to be no obvious pattern in the estimation error and both the estimation error and log estimation error are far from I.I.D.

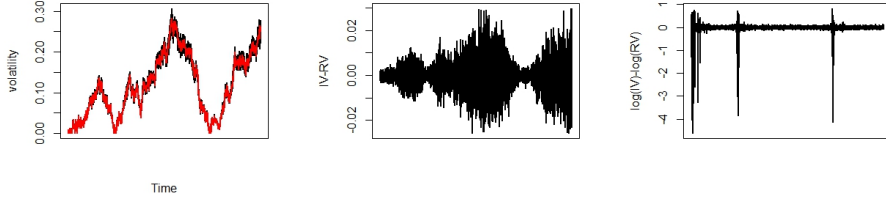


Figure 8: Simulation from model (20). **Left:** The red line represents instantaneous volatility  $\sigma_t$  whereas the black line represents realized volatility  $RV_t$ . **Middle:** Corresponding estimation error for the left simulated path. **Right:** Corresponding log estimation error.

In Figure 9 we have estimated the roughness index via normalized  $p$ -th variation from section 3.1 for realized and instantaneous volatility. In the left plot, we have used the realized volatility as data in the normalized  $p$ -th variation statistic and plotted  $\log(W(K = 500, L = 500 \times 500, \pi, p, t = 1, X = RV))$  against  $H = \frac{1}{p}$ . The right graph is a similar plot using instantaneous volatility instead of realized volatility. The obtained roughness estimators are very different for instantaneous and realized volatility. For realized volatility we obtain an estimated roughness of  $\hat{H}_{L=500 \times 500, K=500}(RV) = 0.323$  which is much lower than the estimated roughness index for instantaneous volatility  $\hat{H}_{L=500 \times 500, K=500}(\sigma) =$

0.499. The true roughness of the model is  $H = 0.5$ . Thus, the roughness estimator is quite accurate when using instantaneous volatility which is in accordance with our results from section 3. However, the estimated roughness when using realized volatility is much rougher than the true roughness. As this is a simulation study we do have any measurement errors, and this rougher behaviour of realized volatility purely comes from the estimation error.

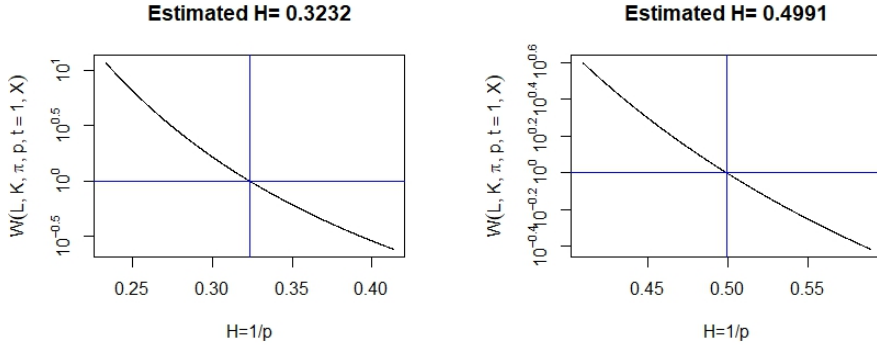


Figure 9: The value of  $\log(W(L = 500 \times 500, K = 500, \pi, p, t = 1, X))$  plotted against  $H = 1/p$  in black. The blue vertical line represents the estimated roughness  $\hat{H}_{L,K}$ . **Left:** Estimated roughness index  $\hat{H}_{L,K}$  for realized volatility derived from the stochastic volatility model (20). **Right:** Estimated roughness index for instantaneous volatility of the same price path.

In Figure 10 we have plotted the estimated roughness index  $\hat{H}_{L=500 \times 500, K}$  against different values of  $K$ . The left graph is for realized volatility while the right graph is for instantaneous volatility. The blue lines represents the estimated roughness  $\hat{H}_{L,K}$  with  $L = 500 \times 500$  and  $K = 500$  which are also reported in Figure 9. From the figure, we observe that irrespective of the choice of  $K$  for the finite sample dataset of length  $L = 500 \times 500$ , the realized volatility is significantly rougher than the instantaneous volatility. The estimated roughness of the volatilities are quite consistent as long as  $K$  is not too low for both volatilities. However,  $\hat{H}_{L=500 \times 500, K}$  seems to be even more consistent for instantaneous volatility than for the realized volatility.

## 4.2 OU-SV model

Now, we will simulate from the following OU-SV model:

$$dS_t = S_t \sigma_t dB_t, \quad \sigma_t = \sigma_0 e^{Y_t}, \quad dY_t = -\gamma Y_t dt + \theta dB'_t, \quad (21)$$

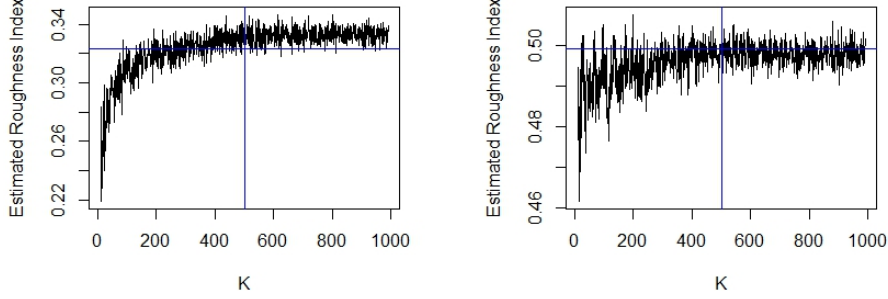


Figure 10: The value of the estimated roughness index  $\hat{H}_{L=500 \times 500, K}$  plotted against different values of  $K$ . **Left:** Estimated roughness index  $\hat{H}_{L=500 \times 500, K}$  for realized volatility derived from the stochastic volatility model (20). The horizontal blue line represents  $\hat{H}_{L=500 \times 500, K=500} = 0.323$ . **Right:** Estimated roughness index for instantaneous volatility of the same price path. The horizontal blue line represents  $\hat{H}_{L=500 \times 500, K=500} = 0.499$ .

where  $S_t$  is the price of the asset and  $B_t$  and  $B'_t$  are two independent Brownian motions. In the simulation, we are using the parameters  $\sigma_0 = \gamma = \theta = 1$  and  $Y_0 = 0$  and we set  $T = 5$  and  $S_0 = 1$ . The true roughness of this model is  $H = 0.5$ .

We are using the same procedure as described in section 4.1 to generate our realized and instantaneous volatility. In Figure 11, we have made plots for these volatilities. The left plot of Figure 11 represents the realized volatility of the price process in black and the instantaneous volatility in red. The middle plot of Figure 11 represents the difference between the realized and instantaneous volatility which is the estimation error. The right plot of Figure 11 represents the log difference between the realized and instantaneous volatility. Visually the middle plot suggests that the estimation error has a complicated dependence structure. However, the log estimation error on the right plot of Figure 11 seems to have an I.I.D. Gaussian structure. This is supported by the theory in Fukasawa et al. [2022]. However, as shown in Section 4.1 the I.I.D. behaviour of the log estimation error does not in general hold for stochastic diffusive models as assumed in Fukasawa et al. [2022].

We now compute the distributions of the estimated roughness index via normalized  $p$ -th statistic  $\hat{H}_{L, K}$  with  $L = 300 \times 300$  and  $K = 300$  for realized volatility and instantaneous volatility. We generate the distributions across 2500 independent paths drawn from the OU-SV model (21). The left plot of Figure 12 is the distribution of  $\hat{H}_{L, K}$  for realized volatility whereas the right plot is the

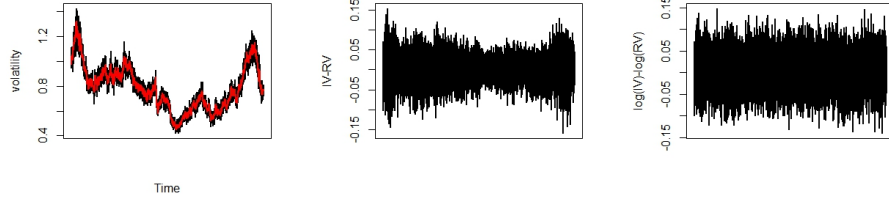


Figure 11: Simulation from the OU-SV model. **Left:** The red line represents instantaneous volatility  $\sigma_t$  whereas the black line represents realized volatility  $RV_t$ . **Middle:** Corresponding estimation error for the left simulated path. **Right:** Corresponding log estimation error.

corresponding distribution for instantaneous volatility as data in the roughness estimator. In addition to this, Table 3 provides the summary statistics for the roughness estimator  $\hat{H}_{L,K}$  for realized and instantaneous volatility respectively across the 2500 independent paths. We observe that realized volatility systematically exhibit much rougher behaviour than instantaneous volatility. The roughness estimates based on instantaneous volatility are very close to the roughness of the true model  $H = 0.5$ . However, the roughness estimates generated from realized volatility are much smaller with a mean of  $\hat{H}_{L,K} = 0.1549$  across the 2500 independent simulations.

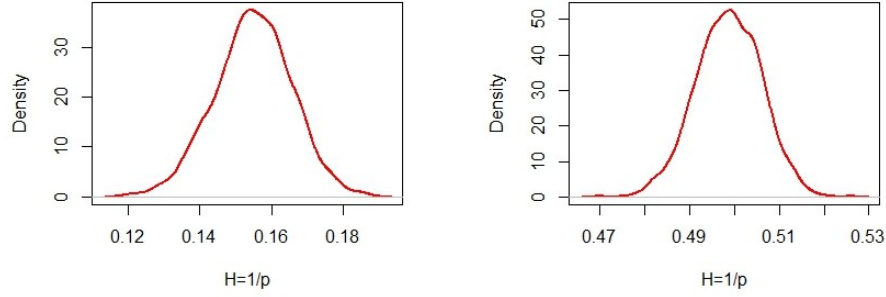


Figure 12: Distribution of the estimated roughness index  $\hat{H}_{L,K}$  via normalized  $p$ -th variation using  $L = 300 \times 300$  and  $K = 300$  across 2500 independent simulations of the OU-SV model (21). **Left:** Realized volatility. **Right:** Instantaneous volatility.

Note that our results deviates slightly from the results in Cont and Das [2024] for a similar model. This might be caused by Cont and Das [2024] using the mean of  $\sigma_t$  across the 300 consecutive data points as their instantaneous volatility which seems to create a slightly smoother volatility process. Furthermore,

	Realized volatility	Instantaneous volatility
Min.	0.1192	0.4698
1st quartile	0.1480	0.4938
Median	0.1551	0.4988
Mean	0.1549	0.4987
3rd quartile	0.1621	0.5039
Max.	0.1873	0.5257

Table 3: Summary of statistics for estimated roughness index  $\hat{H}_{L,K}$  for realized and instantaneous volatility across 2500 independent simulations the OU-SV model (21) with  $L = 300 \times 300$  and  $K = 300$ .

Cont and Das [2024] might be using a different  $T$  than us which in the OU-SV model will significantly impact the size of the estimation error and thus the estimated roughness index for realized volatility when all other parameters are kept constant. The conclusions are however the same. Realized volatility does systematically exhibit much rougher behaviour than instantaneous volatility and than the true roughness of the model. This rougher behaviour is solely caused by the estimation error, and it indicates that roughness estimated based on realized volatility can be unreliable.

### 4.3 A fractional Ornstein-Uhlenbeck model

In the two previous models, instantaneous volatility follows a diffusive behaviour similar to a Brownian motion with  $H = 0.5$ . We now consider a more general case of a stochastic volatility model where the volatility process has a general roughness index  $H \in (0, 1)$  generated from a fractional Brownian motion. We investigate the model for different roughness indexes and explore how this affect the estimation error and estimated roughness index. Consider the following process where the volatility is described by a fractional Ornstein-Uhlenbeck model:

$$dS_t = S_t \sigma_t dB_t, \quad \sigma_t = \sigma_0 e^{Y_t}, \quad dY_t = -\gamma Y_t dt + \theta dB_t^H, \quad (22)$$

where  $S_t$  is the price process,  $B_t$  is a Brownian motion, and  $B_t^H$  is a fractional Brownian motions with Hurst exponent  $H$ . For the simulation we use the parameters  $\sigma_0 = \gamma = \theta = 1$  and  $Y_0 = 0$ , and we set  $T = 1$  and  $S_0 = 1$ .

We now repeat the procedure described in section 4.1 to compute the estimated roughness index  $\hat{H}_{L,K}$  with  $L = 300 \times 300$  and  $K = 300$  for realized and instantaneous volatility. Furthermore, we compute the smoothness parameter from section 3.5 for both the realized and instantaneous volatility. We compute these roughness estimates across different values of the true Hurst exponent  $H$ . In Table (+) we have computed roughness estimates from the fractional Ornstein-Uhlenbeck model (22) with Hurst exponent  $H = \{0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8\}$ .



We have used the same seed in the simulation for all values of  $H$  such that the underlying stochastic element of the price processes are the same. Thus, the difference between the price processes solely lies in the Hurst exponent  $H$ .

## 5 Appendix

### 5.1 A Spectral Method FFT

Consider

$$a_k := \begin{cases} 0 & k = 0; \\ \frac{1}{2} \left( U_{k-1}^{(0)} + iU_{k-1}^{(1)} \right) \sqrt{\frac{f(t_k)}{\ell}} & k = 1, \dots, \ell - 1; \\ U_{k-1}^{(0)} \sqrt{\frac{f(t_k)}{\ell}} & k = \ell; \\ \frac{1}{2} \left( U_{2\ell-k-1}^{(0)} - iU_{2\ell-k-1}^{(1)} \right) \sqrt{\frac{f(t_{2\ell-k})}{\ell}} & k = \ell + 1, \dots, 2\ell - 1. \end{cases}$$

We want to take the Fourier transform of this sequence. The Fourier transform of  $(a_k)_{k=0}^{2\ell-1}$  is  $\lambda_n = \sum_{k=0}^{2\ell-1} a_k \exp(2\pi i \frac{nk}{2\ell})$ . By Euler's formula we obtain

$$\lambda_n = \sum_{k=0}^{2\ell-1} a_k \left( \cos\left(\pi \frac{nk}{\ell}\right) + i \sin\left(\pi \frac{nk}{\ell}\right) \right).$$

We will now insert  $a_k$  and split the sum into the four different cases of  $a_k$ . Hence, we obtain

$$\begin{aligned} \lambda_n &= 0 + \sum_{k=1}^{\ell-1} \frac{1}{2} \left( U_{k-1}^{(0)} + iU_{k-1}^{(1)} \right) \sqrt{\frac{f(t_k)}{\ell}} \left( \cos\left(\pi \frac{nk}{\ell}\right) + i \sin\left(\pi \frac{nk}{\ell}\right) \right) \\ &\quad + U_{\ell-1}^{(0)} \sqrt{\frac{f(t_\ell)}{\ell}} \left( \cos\left(\pi \frac{n\ell}{\ell}\right) + i \sin\left(\pi \frac{n\ell}{\ell}\right) \right) \\ &\quad + \sum_{k=\ell+1}^{2\ell-1} \frac{1}{2} \left( U_{2\ell-k-1}^{(0)} - iU_{2\ell-k-1}^{(1)} \right) \sqrt{\frac{f(t_{2\ell-k})}{\ell}} \left( \cos\left(\pi \frac{nk}{\ell}\right) + i \sin\left(\pi \frac{nk}{\ell}\right) \right). \end{aligned}$$

By definition  $t_k = \frac{\pi k}{\ell}$ . Thus, what is inside the cosine and sine functions is simply  $nt_k$ . Note that for  $k = \ell$  we have  $t_k = \pi$ . Thus,  $\sin(nt_\ell) = 0$  since  $n$  is an integer. By inserting this and changing the indexes of the two sums, we obtain

$$\begin{aligned} \lambda_n &= U_{\ell-1}^{(0)} \sqrt{\frac{f(t_\ell)}{\ell}} \cos(nt_\ell) \\ &\quad + \sum_{k=0}^{\ell-2} \frac{1}{2} \left( U_k^{(0)} + iU_k^{(1)} \right) \sqrt{\frac{f(t_{k+1})}{\ell}} (\cos(nt_{k+1}) + i \sin(nt_{k+1})) \\ &\quad + \sum_{k=0}^{\ell-2} \frac{1}{2} \left( U_{\ell-k-2}^{(0)} - iU_{\ell-k-2}^{(1)} \right) \sqrt{\frac{f(t_{\ell-k-1})}{\ell}} (\cos(nt_{\ell-k-1}) + i \sin(nt_{\ell-k-1})). \end{aligned}$$

Now note that  $\sin(nt_{\ell-k-1}) = -\sin(nt_{\ell-(k+1)})$  since  $t_\ell = \pi$ . Similarly,  $\cos(nt_{\ell-k-1}) = \cos(nt_{\ell-(k+1)})$ . By inserting that in the last sum in the above expression, the

sum becomes

$$\begin{aligned}
& \sum_{k=0}^{l-2} \frac{1}{2} \left( U_{\ell-k-2}^{(0)} - iU_{\ell-k-2}^{(1)} \right) \sqrt{\frac{f(t_{\ell-k-1})}{\ell}} (\cos(nt_{\ell-k-1}) - i \sin(nt_{\ell-k-1})) \\
&= \sum_{k=0}^{l-2} \frac{1}{2} \left( U_k^{(0)} - iU_k^{(1)} \right) \sqrt{\frac{f(t_{k+1})}{\ell}} (\cos(nt_{k+1}) - i \sin(nt_{k+1}))
\end{aligned}$$

where we have reversed the sum order. Now we can combine the two sums into one. Several terms cancel out such that we obtain

$$\begin{aligned}
\lambda_n &= U_{\ell-1}^{(0)} \sqrt{\frac{f(t_\ell)}{\ell}} \cos(nt_\ell) + \sqrt{\frac{f(t_{k+1})}{\ell}} \sum_{k=0}^{l-2} \left( U_k^{(0)} \cos(nt_{k+1}) - U_k^{(1)} \sin(nt_{k+1}) \right) \\
&= \sum_{k=0}^{l-1} \sqrt{\frac{f(t_{k+1})}{\ell}} \left( U_k^{(0)} \cos(nt_{k+1}) - U_k^{(1)} \sin(nt_{k+1}) \right)
\end{aligned}$$

which equals the definition of  $\hat{X}_n^{(\ell)}$  from (11) as desired.

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