# Subsampling realised kernels<sup>\*</sup>

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### Abstract

In a recent paper we have introduced the class of realised kernel estimators of the increments of quadratic variation in the presence of noise. We showed this estimator is consistent and derived its limit distribution under various assumptions on the kernel weights. In this paper we extend our analysis, looking at the class of subsampled realised kernels and we derive the limit theory for this class of estimators. We find that subsampling is highly advantages for estimators based on discontinuous kernels, such as the truncated kernel. For *kinked kernels*, such as the Bartlett kernel, we show that subsampling is impotent, in the sense that subsampling has no effect on the asymptotic distribution. Perhaps surprisingly, for the efficient *smooth kernels*, such as the Parzen kernel, we show that subsampling is harmful as it increases the asymptotic variance. We also study the performance of subsampled realised kernels in simulations and in empirical work.

Keywords: Long run variance estimator; Market frictions; Quadratic variation; Realised kernel; Realised variance; Subsampling.

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# 1 Introduction

High frequency financial data allows us to estimate the increments to quadratic variation, the usual ex-post measure of the variation of asset prices (e.g. Andersen, Bollerslev, Diebold, and Labys (2001) and Barndorff-Nielsen and Shephard (2002)). Common estimators, such as the realised variance, can be sensitive to market frictions when applied to returns recorded over shorter time intervals such as 1 minute, or even more ambitiously, 1 second (e.g. Zhou (1996), Fang (1996) and Andersen, Bollerslev, Diebold, and Labys (2000)). In response two non-parametric generalisations have been proposed: *subsampling* and *realised kernels* by Zhang, Mykland, and Aït-Sahalia (2005) and Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006), respectively. Here we partially unify these approaches by studying the properties of *subsampled realised kernels*.

Our interest is the estimation of the increment to quadratic variation over some arbitrary fixed time period written as [0, t], which could represent a day say, using estimators of the realised kernel type. For a continuous time log-price process X and time gap  $\delta > 0$ , the flat-top<sup>1</sup> realised kernels of Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) take on the following form

$$K(X_{\delta}) = \gamma_0(X_{\delta}) + \sum_{h=1}^{H} k\left(\frac{h-1}{H}\right) \left\{ \gamma_h(X_{\delta}) + \gamma_{-h}(X_{\delta}) \right\}, \quad H \ge 1.$$

Here  $k(x), x \in [0, 1]$ , is a weight function with k(0) = 1, k(1) = 0, while

$$\gamma_h(X_{\delta}) = \sum_{j=1}^{n_{\delta}} x_j x_{j-h}, \quad x_j = X_{\delta j} - X_{\delta(j-1)}, \quad h = -H, ..., -1, 0, 1, ..., H_{\delta}$$

with  $n_{\delta} = \lfloor t/\delta \rfloor$ . Think of  $\delta$  as being small and so  $x_j$  represents the *j*-th high frequency return, while  $\gamma_0(X_{\delta})$  is the realised variance of X. The above authors gave a relatively exhaustive treatment of  $K(X_{\delta})$  when X is a Brownian semimartingale plus noise.

It is important to distinguish three classes of kernels functions k(x): smooth, kinked, and discontinuous. Examples are the Parzen, the Bartlett and the truncated kernel, respectively. Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) have shown that the smooth class, which satisfy k'(0) = k'(1) = 0, lead to realised kernels that converges at the efficient rate,  $n_{\delta}^{1/4}$ . Whereas the kinked kernels, which do not satisfy k'(0) = k'(1) = 0, lead to realised kernels that convergence at  $n_{\delta}^{1/6}$ . The discontinuous kernels lead to inconsistent estimators as we show in Section 3.4.

Realised kernels use returns computed starting at t = 0. There may be efficiency gains by jittering the initial value S times, illustrated in Figure 1, producing S sets of high frequency returns  $x_j^s$ , s = 1, 2, ..., S. Zhang, Mykland, and Aït-Sahalia (2005) made this point for realised variances. We can then average the resulting S realised kernel estimators

$$K(X_{\delta}; S) = \frac{1}{S} \sum_{s=1}^{S} K^{s}(X_{\delta}),$$

<sup>&</sup>lt;sup>1</sup>It is called a flat-top estimator as it imposes that the weight at lag one is one. The motivation for this is discussed extensively in Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006).

where

$$K^{s}(X_{\delta}) = \gamma_{0}^{s}(X_{\delta}) + \sum_{h=1}^{H} k\left(\frac{h-1}{H}\right) \left\{\gamma_{h}^{s}(X_{\delta}) + \gamma_{-h}^{s}(X_{\delta})\right\},$$
  
$$\gamma_{h}^{s}(X_{\delta}) = \sum_{j=1}^{n_{\delta}} x_{j}^{s} x_{j-h}^{s}, \quad x_{j}^{s} = X_{\delta\left(j+\frac{(s-1)}{S}\right)} - X_{\delta\left(j+\frac{(s-1)}{S}-1\right)}$$

We call  $K(X_{\delta}; S)$  the subsampled realised kernel — noting that this form of subsampling is different from the conventional form of subsampling, as we discuss below.

Here we show that subsampling is very useful for the class of discontinuous kernels, because subsampling makes these estimators consistent and converge in distribution at rate  $n^{1/6}$ , where  $n = Sn_{\delta}$  is the *effective sample size*. Zhou (1996) used a simple discontinuous kernel and gave a brief discussion of subsampling that kernel. We will see that his estimator can be made consistent by allowing  $S \to \infty$  as  $n \to \infty$ , a result which is implicit in his paper, but one he did not explicitly draw out. For the class of kinked kernels, we show that subsampling is impotent, in the sense that the asymptotic distribution is the same whether subsampling is used or not. Finally, we show that subsampling is harmful when applied to smooth kernels. In fact, if the number of subsamples, S, increases with the sample size, n, the best rate of convergence is reduced to less than the efficient one,  $n^{1/4}$ .

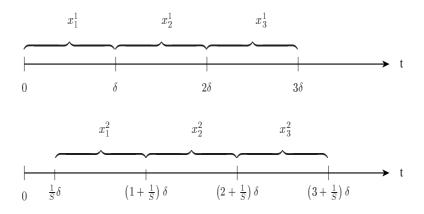


Figure 1:  $x_j^1$  are the usual returns. The bottom series are the offset returns  $x_j^s$ , s = 2, ..., S.

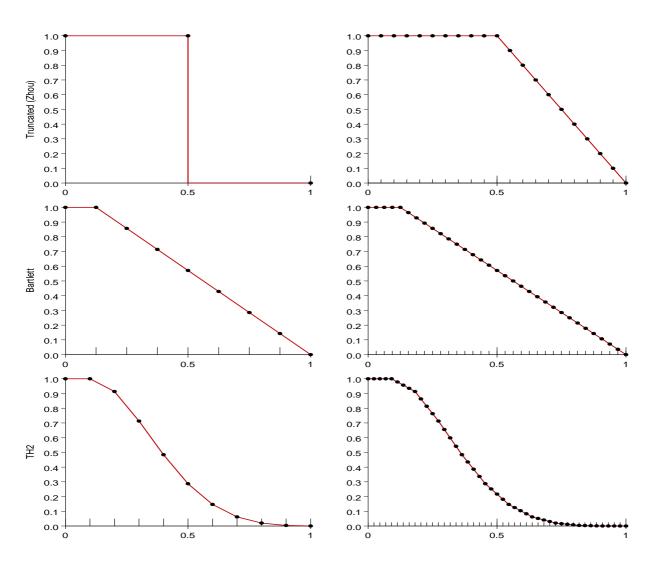
The intuition for these results follows from Lemma A.1 in the appendix. It shows that

$$\gamma_h(X_{\delta}; S) = \frac{1}{S} \sum_{s=1}^{S} \gamma_h^s(X_{\delta}) \simeq \sum_{s=-S+1}^{S-1} k_B\left(\frac{s}{S}\right) \gamma_{Sh+s}(X_{\delta/S}), \quad \text{where} \quad k_B(x) = 1 - |x|,$$

where the approximation is due to subtle end-effects. The implication is that

$$\begin{split} K(X_{\delta};S) &\simeq \sum_{s=-S+1}^{S-1} k_B\left(\frac{s}{S}\right) \gamma_s(X_{\delta/S}) + \sum_{h=1}^H k\left(\frac{h-1}{H}\right) \sum_{s=-S}^S k_B\left(\frac{s}{S}\right) \left\{ \gamma_{Sh+s}(X_{\delta/S}) + \gamma_{-Sh-s}(X_{\delta/S}) \right\} \\ &= \sum_{h=0}^{HS} k_S\left(\frac{h-1}{HS}\right) \tilde{\gamma}_{Sh+s}(X_{\delta/S}). \end{split}$$

So a subsampled realised kernel is a realised kernel simply operating on a higher frequency (ignoring end-effects). The implied kernel weights,  $k_S(\frac{h}{HS})$ ,  $h = 1, \ldots, SH$ , are convex combinations of neighboring weights of the original kernel,



$$k_S\left(\frac{hs}{HS}\right) = \frac{S-s}{S}k\left(\frac{h}{S}\right) + \frac{s}{S}k\left(\frac{h+1}{S}\right), \qquad h = 0, \dots, H, \quad s = 1, \dots, S.$$
(1)

Figure 2: The effects of subsampling some kernels. The left panels display the original kernel functions and the right panels display their implied kernel functions that are induced by subsampling. For the truncated (discontinuous) kernel the two are very different. So subsampling makes an important difference in this case. For the (kinked) Bartlett kernel the two are virtually identical, which suggests that subsampling has no effect on this kernel. Finally, for the smooth kernel in the lower panels subsampling has only a small effect by making the kernel function piecewise linear.

In Figure 2 we trace out the implied kernel weights for three subsampled realised kernels. The left panels display the original kernel functions and right panels display the implied kernel functions. For the truncated kernel (H = 1) subsampling leads to a substantially different implied kernel function – the trapezoidal kernel by Politis and Romano (1995). For the kinked Bartlett kernel subsampling leads to the same kernel function. For a smooth kernel function, the original and implied kernel functions are fairly similar, however subsampling does impose some piecewise linearity which is the reason that subsampling of smooth kernels increases the asymptotic variance.

The connection between subsampled realised kernels and realised kernels is perhaps not too surprising, because Bartlett (1950) motivated his kernel with the subsampling idea. The conventional form of subsampling that is based on subseries that consist of consecutive observations. This is different from our subsamples that consist of every Sth observation. Such are called subgrids in Zhang, Mykland, and Aït-Sahalia (2005). While the two types of subsampling are different they can result in identical estimators. For instance, the sparsely sampled realised variance,  $\gamma_0^1(X_\delta)$ , is identical to Carlstein's subsample estimator (of the variance of a sample mean when the mean is zero), see Carlstein (1986). Carlstein's estimator is based on non-overlapping subseries and Künsch (1989) analysed the closely related estimator based on overlapping subseries. Interestingly, the (overlapping) subsample estimator by Künsch (1989) is identical to the average sparsely sampled realised variance called "second best" in Zhang, Mykland, and Aït-Sahalia (2005), so that the TSRV and MSRV estimators, by Zhang, Mykland, and Aït-Sahalia (2005), Aït-Sahalia, Mykland, and Zhang (2006), and Zhang (2006), can be expressed as linear combinations of two or more subsample estimators of the overlapping subseries type by Künsch (1989). For additional details on the relation between Bartlett kernel and various subsample estimators, see Anderson (1971, p. 512), Priestley (1981, pp. 439–440), and Politis, Romano, and Wolf (1999, pp. 95–98).

This paper has the following structure. We present the basic framework in Section 2 along with some known results. In Section 3 we present our main results. Here we derive the limit theory for subsampled realised kernels and show that subsampling cannot improve realised kernels within a very broad class of estimators. In Section 4, we given some specific recommendations on empirical implementation of subsampled realised kernels and how to conduct valid inference in this context. We present results from a small simulation study in Section 5 and an empirical application in Section 6. We conclude in Section 7 and present all proofs in an appendix.

# 2 Notation, definitions and background

### 2.1 Semimartingales and quadratic variation

The fundamental theory of asset prices says that the log-price at time  $t, Y_t$ , must, in a frictionless arbitrage free market, obey a *semimartingale* process (written  $Y \in SM$ ) on some filtered probability space  $(\Omega, \mathcal{F}, (\mathcal{F}_t)_{t \geq T^*}, P)$ , where  $T^* \leq 0$ . Crucial to semimartingales, and to the economics of financial risk, is the quadratic variation (QV) process of  $Y \in SM$ . This can be defined as

$$[Y]_{t} = \lim_{N \to \infty} \sum_{j=1}^{N} \left( Y_{t_{j}} - Y_{t_{j-1}} \right)^{2}, \qquad (2)$$

(e.g. Protter (2004, p. 66–77) and Jacod and Shiryaev (2003, p. 51)) for any sequence of deterministic partitions  $0 = t_0 < t_1 < ... < t_N = t$  with  $\sup_j \{t_{j+1} - t_j\} \to 0$  for  $N \to \infty$ .

The most familiar semimartingales are of Brownian semimartingale type  $(Y \in \mathcal{BSM})$ 

$$Y_t = \int_0^t a_u \mathrm{d}u + \int_0^t \sigma_u \mathrm{d}W_u,\tag{3}$$

where a is a predictable locally bounded drift,  $\sigma$  is a càdlàg volatility process and W is a Brownian motion. If  $Y \in \mathcal{BSM}$  then  $[Y]_t = \int_0^t \sigma_u^2 du$ . In some of our asymptotic theory we also assume, for simplicity of exposition, that

$$\sigma_t = \sigma_0 + \int_0^t a_u^{\#} du + \int_0^t \sigma_u^{\#} dW_u + \int_0^t v_u^{\#} dV_u,$$
(4)

where  $a^{\#}$ ,  $\sigma^{\#}$  and  $v^{\#}$  are adapted càdlàg processes, with  $a^{\#}$  also being predictable and locally bounded and V is Brownian motion independent of W. Much of what we do here can be extended to allow for jumps in  $\sigma$  (cf. Barndorff-Nielsen, Graversen, Jacod, and Shephard (2006)).

### 2.2 Assumptions about noise

We write the effects of market frictions as U, so that we observe the process

$$X = Y + U. (5)$$

Our scientific interest will be in estimating  $[Y]_t$ . In the main part of our work we will assume that  $Y \perp U$  where, in general,  $A \perp B$  denotes that A and B are independent. From a market microstructure theory viewpoint this is a strong assumption as one may expect U to be correlated with increments in Y. However, the empirical work of Hansen and Lunde (2006) suggests this independence assumption is not too damaging statistically when we analyse data in thickly traded stocks recorded approximately every minute (see also Kalnina and Linton (2006)).

We make a white noise assumption about the U process  $(U \in \mathcal{WN})$ :

$$\mathbf{E}(U_t) = 0, \quad \operatorname{Var}(U_t) = \omega^2, \quad \operatorname{Var}(U_t^2) = \lambda^2 \omega^4, \quad U_t \perp \!\!\!\perp U_s \tag{6}$$

for any  $t \neq s$ , where  $\lambda \in \mathbb{R}^+$ . This white noise assumption is unsatisfactory but is a useful starting point if we think of the market frictions as operating in tick time (e.g. Bandi and Russell (2005), Zhang, Mykland, and Aït-Sahalia (2005) and Hansen and Lunde (2006)).

Analogous to the realised autocovariances we define

$$\gamma_h(Y_{\delta}, U_{\delta}) = \sum_{j=1}^{n_{\delta}} y_j u_{j-h}, \quad y_j = Y_{\delta j} - Y_{\delta(j-1)} \text{ and } u_j = U_{\delta j} - U_{\delta(j-1)}.$$

From (5) we have that

$$\gamma_h(X_{\delta}) = \gamma_h(Y_{\delta}) + \gamma_h(Y_{\delta}, U_{\delta}) + \gamma_h(U_{\delta}, Y_{\delta}) + \gamma_h(U_{\delta})$$

It will be useful to have the following notation  $\widetilde{\gamma}(X_{\delta}) = \{\gamma_0(X_{\delta}), \widetilde{\gamma}_1(X_{\delta}), ..., \widetilde{\gamma}_H(X_{\delta})\}^{\mathsf{T}}$ , where  $\widetilde{\gamma}_h(X_{\delta}) = \gamma_h(X_{\delta}) + \gamma_{-h}(X_{\delta})$ , and introduce the analogous definitions of  $\widetilde{\gamma}(Y_{\delta}), \widetilde{\gamma}(U_{\delta})$ , and  $\widetilde{\gamma}(Y_{\delta}, U_{\delta})$ .

# 3 Subsampled realised kernel

Here we study subsampled realised kernels based on smooth and kinked kernel functions. Specifically, we require that k(s) is continuous and twice differentiable on [0, 1] and that k(0) = 1 and k(1) = 0. Naturally, the derivatives at the end points are defined by  $k'(0) = \lim_{x \downarrow 0} \frac{k(x) - k(0)}{x}$  and  $k'(1) = \lim_{x \uparrow 1} \frac{k(1) - k(x)}{1 - x}$ .

Without subsampling, Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) showed that

$$k'(0) = 0$$
 and  $k'(1) = 0,$  (7)

is a necessary condition for a realised kernel to have the best rate of convergence, and this property is also key for subsampled realised kernels — see also the work of Zhang (2006) on using subsampling of realised variance to obtain the same rate of convergence. We shall refer to continuous kernels that satisfy (7) as *smooth*, otherwise they are called *kinked*.

In some of our proofs it is convenient to extend the support of the kernel functions beyond the unit interval, using the conventions: k(x) = 0 for x > 1 and k(-x) = k(x).

Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) showed that kernel functions of the type can be used to produce consistent estimators with mixed Gaussian asymptotic distributions. It is therefore interesting to analyze whether there are any gain from subsampling realised kernels or not. Perhaps surprisingly we find that subsampling is harmful or, at best, impotent, for realised kernels that are based on smooth or kinked kernel functions.

Below we formulate limit results for subsampled realised kernels using the notation

$$k_{\bullet}^{0,0} = \int_{0}^{1} k(x)^{2} dx, \quad k_{\bullet}^{1,1} = \int_{0}^{1} k'(x)^{2} dx, \quad k_{\bullet}^{2,2} = \int_{0}^{1} k''(x)^{2} dx,$$
  

$$\xi = \omega^{2} / \sqrt{t \int_{0}^{t} \sigma_{u}^{4} du}, \quad \rho = \int_{0}^{t} \sigma_{u}^{2} du / \sqrt{t \int_{0}^{t} \sigma_{u}^{4} du},$$

and we define  $\tilde{K}(X_{\delta}; S) = K(X_{\delta}; S) + \Delta_{H,n}^{S}$ , where  $\Delta_{H,n}^{S} = S^{-1} \sum_{s=1}^{S} \sum_{h=1}^{H} \left\{ k(\frac{h+1}{H}) - k(\frac{h-1}{H}) \right\} R_{h,n}^{s}$ ,  $R_{h,n}^{s} = \frac{1}{2} (U_{t_{n}^{s}} U_{t_{n}^{s} + h\delta} + U_{t_{0}^{s}} U_{t_{0}^{s} - h\delta} - U_{t_{n}^{s}} U_{t_{n}^{s} - h\delta} - U_{t_{0}^{s}} U_{t_{0}^{s} + h\delta}), t_{0}^{s} = \frac{s-1}{S} \delta$ , and  $t_{n}^{s} = t + \frac{s-1}{S} \delta$ . So  $\Delta_{H,n}^{S}$  is related to end-effects.

**Theorem 1** For large H and n the asymptotic distributions of

$$K(Y_{\delta};S) - \int_0^t \sigma_u^2 \mathrm{d}u, \quad K(Y_{\delta},U_{\delta};S) + K(U_{\delta},Y_{\delta};S), \quad and \quad K(U_{\delta};S) + \Delta_{H,n}^S,$$

are mixed Gaussian, uncorrelated with mean zero and asymptotic variances given by

$$4\frac{H}{n_{\delta}}k_{\bullet}^{0,0}t\int_{0}^{t}\sigma_{u}^{4}\mathrm{d}u,\tag{8}$$

$$8\omega^2 \int_0^t \sigma_u^2 \mathrm{d}u k_{\bullet}^{1,1} H^{-1} \middle/ S \tag{9}$$

$$4\omega^4 n_\delta \left[ \left\{ k'(0)^2 + k'(1)^2 \right\} H^{-2} + k_{\bullet}^{2,2} H^{-3} \right] / S.$$
(10)

respectively, and the asymptotic variance of  $\Delta_{H,n}^S$  is  $4\omega^4 k_{\bullet}^{1,1}/(HS)$ . Furthermore,  $\tilde{K}(X_{\delta};S) - \int_0^t \sigma_u^2 du$  is mixed Gaussian with a zero mean and variance

$$4t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \left\{ \frac{H}{n_{\delta}} k_{\bullet}^{0,0} + \frac{2\xi \rho k_{\bullet}^{1,1} H^{-1} + \xi^{2} n_{\delta} \left[ \left\{ k'(0)^{2} + k'(1)^{2} \right\} H^{-2} + k_{\bullet}^{2,2} H^{-3} \right]}{S} \right\}.$$
(11)

Subsampling has no impact on the first term, (8). This is despite the fact that subsampling lowers the variance of the individual realised autocovariances,  $\tilde{\gamma}_h(Y_\delta)$ . This is because subsampling introduces positive correlation between  $\tilde{\gamma}_h(Y_\delta; S)$  and  $\tilde{\gamma}_{h+1}(Y_\delta; S)$  that exactly offsets the reduction in the variance of the realised autocovariances. Subsampling does reduce the variances of the terms effected by noise, (9) and (10), by a factor of S.

The auxiliary quantity,  $\tilde{K}(X_{\delta}; S)$ , is introduced to simplify the exposition of our results.  $\tilde{K}(X_{\delta}; S)$ and  $K(X_{\delta}; S)$  are often asymptotically equivalent because their difference,  $\Delta_{H,n}^{S}$ , vanishes at a sufficiently fast rate. This is made precise in the following Lemma.

**Lemma 1** If 
$$k'(0)^2 + k'(1)^2 \neq 0$$
 or  $S \to \infty$ , then  $\operatorname{avar}\{K(X_{\delta}) - \tilde{K}(X_{\delta})\}/\operatorname{avar}\{\tilde{K}(X_{\delta})\} = o(1)$ . If  $k'(0)^2 + k'(1)^2 = 0$  then  $\operatorname{avar}\{K(X_{\delta}) - \tilde{K}(X_{\delta})\}/\operatorname{avar}\{\tilde{K}(X_{\delta})\} \leq \xi/\left\{2 + 2\sqrt{k_{\bullet}^{2,2}k_{\bullet}^{0,0}/(k_{\bullet}^{1,1})^2}\right\}$ .

We shall state several asymptotic results for  $n^{\gamma} \left\{ \tilde{K}(X_{\delta}) - \int_{0}^{t} \sigma_{u}^{2} du \right\}$ . An implication of Lemma 1 is that  $K(X_{\delta})$  can be substituted for  $\tilde{K}(X_{\delta})$  whenever  $\gamma < 1/4$ . When  $\gamma = 1/4$  the difference between  $K(X_{\delta})$  and  $\tilde{K}(X_{\delta})$  is not trivial in an asymptotic sense, but for all practical purposes their difference is negligible. The reason being that a realistic empirical value for  $\xi$ , is  $\xi \leq 0.01$ . With the original Tukey-Hanning kernel the relative variance in Lemma 1 is no larger than  $1/\{200(1+\sqrt{3})\} \approx 0.00183$ .

The most obvious generalisation of Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) is to think of the case where S is fixed and we allow H to increase with  $n_{\delta}$ . When (7) holds, we can follow Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) and set  $H = c(\xi n_{\delta})^{1/2}$ . Then we obtain the result that, where Ls denotes convergence in law stably (e.g. Barndorff-Nielsen, Graversen, Jacod, and Shephard (2006)),

$$n_{\delta}^{1/4} \left\{ \tilde{K}(X_{\delta}; S) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left\{ 0, 4\omega \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{3/4} \left( ck_{\bullet}^{0,0} + \frac{2c^{-1}\rho k_{\bullet}^{1,1} + c^{-3}k_{\bullet}^{2,2}}{S} \right) \right\}.$$

Whether or not (7) holds, when we set  $H = c(\xi n_{\delta})^{2/3}$  we have

$$n_{\delta}^{1/6} \left\{ \tilde{K}(X_{\delta}; S) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left[ 0, 4\omega^{4/3} \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{2/3} \left\{ ck_{\bullet}^{0,0} + \frac{k'(0)^{2} + k'(1)^{2}}{c^{2}S} \right\} \right].$$

Here S plays a relatively simple role, reducing the impact of noise — by in effect reducing the noise variance from  $\omega^2$  to  $\omega^2/\sqrt{S}$ . If (7) does hold then we get

$$n_{\delta}^{1/6} \left\{ \tilde{K}(X_{\delta}; S) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left\{ 0, 4ck_{\bullet}^{0,0} \omega^{4/3} \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{2/3} \right\},$$

which implies no asymptotic gains at all from subsampling.

### 3.1 Effective Sample Size

The effectiveness of subsampling can be assessed in terms of the effective sample size,  $n = n_{\delta}S$ . It makes explicit that a larger S reduces the sample size,  $n_{\delta}$ , that is available for each of the realised kernels. Then we ask if it is better to increase  $n_{\delta}$  or S for a given n. In terms of n (11) becomes

$$4t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \left[ \frac{HS}{n} k_{\bullet}^{0,0} + \frac{2\xi \rho k_{\bullet}^{1,1}}{HS} + n\xi^{2} \left\{ \frac{k'(0)^{2} + k'(1)^{2}}{(HS)^{2}} + S \frac{k_{\bullet}^{2,2}}{(HS)^{3}} \right\} \right].$$
(12)

Here HS appears in the variance expression in a way that is almost identical to H when there is no subsampling (S = 1). The only difference is the impact on the last term. This term vanishes when k'(0) = k'(1) = 0 does not hold, because the second last term is then  $O(n/(SH)^2)$  whereas the last term is only  $O(H^{-1}) O(n/(SH)^2)$ . This feature of the asymptotic variance holds the key to the different results we derive for smooth and kinked kernels.

### **3.2** Kinked Kernels: When k'(0) = k'(1) = 0 does not hold

When (7) does not hold the asymptotic variance of  $K(X_{\delta}, S)$  is given by

$$4t \int_0^t \sigma_u^4 \mathrm{d}u \left\{ \frac{HS}{n} k_{\bullet}^{0,0} + \frac{2\xi\rho k_{\bullet}^{1,1}}{HS} + n\xi^2 \frac{k'(0)^2 + k'(1)^2}{(HS)^2} \right\}.$$

While this expression depends on the product HS, it is invariant to the particular values of H and S, though we do need  $H \to \infty$  to justify the terms,  $k_{\bullet}^{0,0}$ ,  $k_{\bullet}^{1,1}$ , etc. We have the following result.

**Theorem 2** (i) If  $SH = c(\xi n)^{2/3}$  we have

$$n^{1/6} \left( \tilde{K}(X_{\delta}; S) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right) \xrightarrow{L_{s}} MN \left( 0, 4\omega^{4/3} \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{2/3} \left\{ ck_{\bullet}^{0,0} + \frac{k'(0)^{2} + k'(1)^{2}}{c^{2}} \right\} \right), \quad (13)$$

as  $n \to \infty$ , so long as H increase with n. (ii) The asymptotic variance is minimised by

$$c = \left\{ 2 \frac{k'(0)^2 + k'(1)^2}{k_{\bullet}^{0,0}} \right\}^{1/3}, \quad and \quad 6ck_{\bullet}^{0,0}\omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3}$$

is the lower bound for the asymptotic variance.

Thus (13) is not influenced by S, not even the rate of growth in S. All that matters is that H grows and that HS grows at the right rate. The implication is that there are no gains from subsampling when  $k'(0)^2 + k'(1)^2 \neq 0$ . So we might as well set S = 1 and use the realised kernel that does not require any subsampling. The second part of Theorem 2 shows that

$$ck_{\bullet}^{0,0} = 6 \left[ 2 \left( k_{\bullet}^{0,0} \right)^2 \left\{ k'(0)^2 + k'(1)^2 \right\} \right]^{1/3}$$

controls the asymptotic efficiency of estimators in this class.

**Example 1** The Bartlett kernel, k(x) = 1 - x, has  $k_{\bullet}^{0,0} = 1/3$  and  $k'(0)^2 + k'(1)^2 = 2$ , so that  $6ck_{\bullet}^{0,0} = 2 \cdot 12^{1/3} \simeq 4.58$ , whereas the quadratic kernel,  $k(x) = 1 - 2x + x^2$ , is more efficient, because it has  $k_{\bullet}^{0,0} = 1/5$  and  $k'(0)^2 + k'(1)^2 = 4$ , so that  $6ck_{\bullet}^{0,0} = 12 \cdot 5^{-2/3} \simeq 4.10$ .

# **3.3 Smooth Kernels: When** k'(0) = k'(1) = 0 holds

In this Section we consider smooth kernel functions. Some examples of smooth kernel functions are given in Table 1, where  $k_{\text{TH}_1}(x)$  is the Tukey-Hanning kernel.

Cubic kernel	$k_{C}(x)=1-3x^{2}+2x^{3}$
Parzen kernel	$k_P(x) = \begin{cases} 1 - 6x^2 + 6x^3 & 0 \le x \le 1/2\\ 2(1-x)^3 & 1/2 \le x \le 1 \end{cases}$
$\mathrm{TH}_p$	$k_{TH_p}(x) = \sin^2 \{ \pi/2 (1-x)^p \}$

Table 1: Some smooth kernel functions.

We know from Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) that the rate of convergence of realised kernels improves when k'(0) = k'(1) = 0. This smoothness condition will also improve the rate of convergence for subsampled realised kernels. For smooth kernel functions, the asymptotic variance is given by

$$4t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \left\{ \frac{HS}{n} k_{\bullet}^{0,0} + \frac{2\xi \rho k_{\bullet}^{1,1}}{HS} + \xi^{2} nS \frac{k_{\bullet}^{2,2}}{(HS)^{3}} \right\}.$$
 (14)

Because the last term is multiplied with S it is evident that the asymptotic distribution will depend on whether S is constant or increases with n. This is made precise in the following Theorem.

**Theorem 3** (i.a) When S is fixed we set  $HS = c(\xi n)^{1/2}$  and have

$$n^{1/4} \left\{ \tilde{K}(X_{\delta}) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left[ 0, 4\omega \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{3/4} \left\{ ck_{\bullet}^{0,0} + \frac{2\rho}{c}k_{\bullet}^{1,1} + \frac{S}{c^{3}}k_{\bullet}^{2,2} \right\} \right].$$
(15)

(i.b) When  $S = an^{\alpha}$  for some  $0 < \alpha < 2/3$ , we set  $HS = c(\xi n)^{1/2}n^{\alpha/4}$  and have

$$n^{\frac{1-\alpha/2}{4}} \left( \tilde{K}(X_{\delta}; S) - \int_{0}^{t} \sigma_{u}^{2} \mathrm{d}u \right) \xrightarrow{L_{\delta}} MN \left[ 0, 4\omega \left( t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \right)^{3/4} \left\{ ck_{\bullet}^{0,0} + \frac{a}{c^{3}}k_{\bullet}^{2,2} \right\} \right]$$

(ii) Whether S is constant or not, the asymptotic variance is minimized by

$$HS = (\xi n)^{1/2} \sqrt{\frac{\rho k_{\bullet}^{1,1}}{k_{\bullet}^{0,0}}} \left\{ 1 + \sqrt{1 + 3S \frac{k_{\bullet}^{0,0} k_{\bullet}^{2,2}}{(\rho k_{\bullet}^{1,1})^2}} \right\}$$

and the lower bound is

$$n^{-1/2}\omega\left(t\int_0^t\sigma_u^4\mathrm{d}u\right)^{3/4}g(S),\tag{16}$$

where

$$g(S) = \frac{16}{3} \sqrt{\rho k_{\bullet}^{1,1} k_{\bullet}^{0,0}} \left\{ \frac{1}{\sqrt{1 + \sqrt{1 + 3S \frac{k_{\bullet}^{0,0} k_{\bullet}^{2,2}}{(\rho k_{\bullet}^{1,1})^2}}} + \sqrt{1 + \sqrt{1 + 3S \frac{k_{\bullet}^{0,0} k_{\bullet}^{2,2}}{(\rho k_{\bullet}^{1,1})^2}} \right\}.$$
 (17)

**Remark.** In (*i.b*) we impose  $\alpha < 2/3$ . The reason is that  $H \propto n^{1/2 + \alpha/4 - \alpha} = n^{(1 - \frac{3}{2}\alpha)/2}$  and we need  $(1 - \frac{3}{2}\alpha)/2 > 0$  to ensure that H grows with n.

The relative efficiency in this class of estimators is given from g(S), and we have the following important result for subsampling of smooth kernels

### **Corollary 1** The asymptotic variance of $\tilde{K}(X_{\delta}; S)$ is strictly increasing in S.

The implication is that subsampling is always harmful for smooth kernels. Furthermore, (i.b) shows that there is an efficiency loss from allowing S to grow with n. See Table 2 for the values of g(S) for some selected kernel functions.

Another implication of Theorem 3 concerns the best way to sample high frequency returns. This result is formulated in the next corollary and will require some explanation.

### **Corollary 2** The asymptotic variance, (16), as a function of $\rho$ , is minimized for $\rho = 1$ .

The Corollary is interesting because  $\rho = \int_0^t \sigma_u^2 du / \sqrt{t \int_0^t \sigma^4 du}$  depends on the sampling scheme by which intraday returns are obtained. So  $\rho$  can be interpreted as an asymptotic measure of heteroskedasticity in the intraday returns, where  $\rho = 1$  corresponds to homoskedastic intraday returns. Rather than equidistant sampling in calendar time we can generate the sampling times by,

$$t_j = t \times \tau\left(\frac{j}{n}\right), \qquad j = 0, 1, \dots, n,$$

where  $\tau$  is a time change  $(\tau(0) = 0, \tau(1) = 1, \text{ and } \tau$  is monotonically increasing, so that  $0 = t_0 \leq t_1 \leq \cdots \leq t_n = t$ ). A change of time does not affect  $\int_0^t \sigma_u^2 du$  but does influence the integrated quarticity  $\int_0^t \sigma_u^4 du$ , see e.g. Mykland and Zhang (2006). A particularly interesting sampling scheme is business time sampling (BTS), see e.g. Oomen (2005, 2006), which is the sampling scheme that minimises the integrated quarticity, see Hansen and Lunde (2006, p. 135). It is easy to see that the time change associated with BTS,  $\tau(\cdot)$ ,  $\tau_{\text{BTS}}(\cdot)$  say, must solve  $\int_0^{t\times\tau(s)} \sigma_u^2 du = s \times \int_0^t \sigma_u^2 du$ , and by the *implicit function theorem* we have  $\tau'_{\text{BTS}}(s) \propto 1/\sigma^2(\tilde{s})$ , where  $\tilde{s} = t \times \tau_{\text{BTS}}(s)$ . The implication is that returns are sampled more frequently when the volatility is high and less frequently when the volatility is low under BTS. In general we have  $\rho \leq 1$  and Corollary 2 shows that BTS ( $\rho = 1$ ) is the ideal sample scheme. Naturally, sampling in business time is infeasible because  $\tau_{\text{BTS}}$  depends on the unknown volatility path. Still, Corollary 2 can be used as argument in favor of sampling schemes that results in less heteroskedastic intraday returns than does CTS.

Given S and  $\rho$  the optimal H is  $H = c_S(\xi n)^{1/2}$  for this class of kernels where

$$c_S = S^{-1} \sqrt{\frac{\rho k_{\bullet,\bullet}^{1,1}}{k_{\bullet,\bullet}^{0,0}}} \left\{ 1 + \sqrt{1 + 3S \frac{k_{\bullet,\bullet}^{0,0} k_{\bullet,\bullet}^{2,2}}{(\rho k_{\bullet,\bullet}^{1,1})^2}} \right\}.$$
 (18)

In Table 2 we present key quantities for some smooth kernels. Perhaps the most interesting quantity is g(S) of (17), as it enable us to compare the relative efficiency across estimators. In Table 2 we have computed g(S) for the case where  $\rho = 1$ . So g(S) can be compared to 8 which

	$k^{0,0}_{ullet}$	$k_{\bullet}^{1,1}$	$k_{ullet}^{2,2}$	$\sqrt{k_{ullet}^{0,0}k_{ullet}^{1,1}}$	$\frac{k_{\bullet}^{0,0}k_{\bullet}^{2,2}}{(k_{\bullet}^{1,1})^2}$	$c_1$	g(S)				
					(~•)		S = 1	S = 2	S = 3	S = 10	
Cubic	0.371	1.20	12.0	0.67	3.09	3.68	9.03	9.81	10.39	12.72	
Parzen	0.269	1.50	24.0	0.64	2.87	4.77	8.53	9.25	9.78	11.94	
$\mathrm{TH}_1$	0.375	1.23	12.2	0.68	3.00	3.70	9.18	9.96	10.55	12.89	
$\mathrm{TH}_2$	0.218	1.71	41.8	0.61	3.11	5.75	8.27	8.99	9.51	11.65	
${ m TH}_5$	0.097	3.50	489.0	0.58	3.85	8.07	8.07	8.82	10.19	11.57	
$\mathrm{TH}_{10}$	0.050	6.57	3610.6	0.57	4.19	24.79	8.04	8.80	10.19	11.59	
${ m TH_{16}}$	0.032	10.26	14374.0	0.57	4.33	39.16	8.02	8.80	10.20	11.60	

Table 2: Key quantities for some smooth-continuous kernels.

Key is g(S) that measures the relative efficiency in this class of estimators. Here computed for the case with constant volatility ( $\rho = 1$ ) such that these numbers are comparable with the maximum likelihood estimator that has g = 8.00. No subsampling (S = 1) produces the best estimator and kernels with a relative large  $k_{\bullet}^{0,0} k_{\bullet}^{2,2} / (k_{\bullet}^{1,1})^2$  tend to be more sensitive to subsampling.

is the corresponding constant for the maximum likelihood estimator in the Gaussian parametric version of the problem. We see that most kernels are only slightly less efficient than the maximum likelihood estimator, TH<sub>16</sub> almost reaching this lower bound. Comparing g(S) for different degrees of subsampling, reminds us that S = 1 (no subsampling) yields the most efficient estimator. The larger the value of  $k_{\bullet}^{0,0}k_{\bullet}^{2,2}/(k_{\bullet}^{1,1})^2$  the more sensitive is the kernel to subsampling.

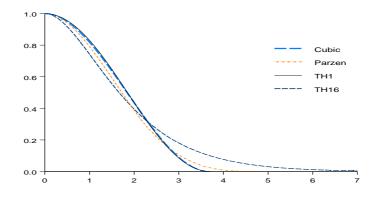


Figure 3: Some smooth kernels,  $k(x/c_1)$ , using their respective optimal value of c when S = 1.

Figure 3 plots some smooth kernel functions,  $k(x/c_1)$  using their respective optimal value for  $c_1$ , see Table 2. We see that the TH<sub>1</sub> kernel is almost identical to the cubic kernel. The TH<sub>16</sub> kernel is somewhat flatter, putting less weight on realised autocovariances of lower order and higher weight on realised autocovariances of higher order. The Parzen kernel is typically between TH<sub>1</sub> and TH<sub>16</sub>.

While the smooth kernels improve the rate of convergence over the kinked kernels, the improvements may be modest in finite samples. The reason is the following. When the noise is small the optimal H is small, and H may actually be quite similar for kinked and smooth kernels. For instance with  $\xi = 0.01$  and n = 780, the Bartlett kernel has  $c_{\text{BARTLETT}}(\xi n)^{2/3} = 9.00$  whereas the cubic kernel has  $c_{\text{CUBIC}}(\xi n)^{1/2} = 10.78$ . So in this case the two types of estimators are rather similar and despite the fact that  $H_{\text{BARTLETT}}$  grows at the faster rate  $n^{2/3}$ , the cubic kernels includes more lags in this situation.

#### **3.4** Discontinuous kernel functions

In this section we consider the kernel functions we have labelled as discontinuous kernels. Such kernels lead to estimators with poor asymptotic properties. We shall see that subsampling can substantially improve such estimators, making them consistent with mixed Gaussian distributions. So for such kernels, subsampling is a saviour.

**Lemma 2** Let  $K_w(X_{\delta}) = \sum_{h=0}^{H} w_h \tilde{\gamma}_h(X_{\delta})$ , where H = o(n) (possibly constant). Then  $w_0 = 1 + o(1)$  and  $w_0 - w_1 = o(n^{-1})$ , are necessary conditions for  $E(K_w(X_{\delta}) - \int_0^t \sigma_u^2 du) \to 0$ ; and

$$\sum_{h=0}^{H} (w_{h+1} - 2w_h + w_{h-1})^2 = o(n^{-1}),$$
(19)

is a necessary condition for  $\operatorname{Var}\left(K_w(X_{\delta}) - \int_0^t \sigma_u^2 \mathrm{d}u\right) \to 0$ , where we set  $w_{H+1} = 0$  and  $w_{-1} = w_0$ .

The lemma shows that realised kernels using a fixed H cannot converge to  $\int_0^t \sigma_u^2 du$  in mean squares, because such estimators will not satisfy (19).

Consider the case where we construct  $w_h$  from a kernel function and let  $H \to \infty$ . In this situation it is clear that any discontinuous kernel will violate (19), because

$$n\sum_{h=0}^{H} (w_{h+1} - 2w_h + w_{h-1})^2 \simeq n \times \sum_{x_j \in \mathcal{D}_k} \left\{ \lim_{x \uparrow x_j} k(x) - \lim_{x \downarrow x_j} k(x) \right\}^2.$$

Here  $\mathcal{D}_k$  is the set of discontinuity points for k(x).

Next, we consider the truncated kernel which does not satisfies (19). We will see that subsampling this kernel produces an estimator that is consistent and mixed Gaussian. This is true whether H is finite or is allowed to grow with the sample size.

### **3.4.1** Zhou (1996) estimator

First we will look at estimators which are thought of as having H fixed and allowing the degree of subsampling to increase. This is outside the spirit of the realised kernels of Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) which need H to get large with  $n_{\delta}$  for consistency, however it is close to the important early work of Zhou (1996) and is strongly intellectually connected to the two scale estimators by Zhang, Mykland, and Aït-Sahalia (2005).

The Zhou (1996) estimator can be written as  $\gamma_0(X_{\delta}; S) + \tilde{\gamma}_1(X_{\delta}; S)$  which is the subsampled realised kernel based on the truncated kernel function using H = 1. Zhou (1996) noticed that the variance of his estimator was of order  $O(\frac{S}{n_{\delta}}) + O(\frac{1}{S}) + O(\frac{n_{\delta}}{S^2})$ , but did not realize that by allowing S to increase with  $n_{\delta}$  his estimator is consistent. In fact, in a subsequent paper Zhou stated that his subsampled realised kernels was inconsistent, see Zhou (1998, p. 114). The following Theorem gives its asymptotic distribution. **Theorem 4** Suppose  $S = c^3 n_{\delta}^2$ , then as  $n_{\delta} \to \infty$ 

$$n_{\delta}^{1/2} \left\{ \gamma_0(X_{\delta}; S) + \widetilde{\gamma}_1(X_{\delta}; S) - \int_0^t \sigma_u^2 \mathrm{d}u \right\} \xrightarrow{Ls} MN\left(0, \frac{16}{3}t \int_0^t \sigma_u^4 \mathrm{d}u + 8\omega^4/c^3\right).$$

This asymptotics is not particularly attractive for its seeming  $n_{\delta}^{1/2}$  rate of convergence hides the fact that it assumes massive databases in order to allow S to increase rapidly with  $n_{\delta}$ . A more interesting way of thinking about this estimator is in terms of the effective sample size  $n = S \times n_{\delta}$ . Again we ask if it is better to increase  $n_{\delta}$  or S for a given n. This leads to the following result.

**Lemma 3** If  $S = c(\xi n)^{2/3}$  then the Zhou estimator has

$$n^{1/6} \left( \gamma_0(X_{\delta}; S) + \widetilde{\gamma}_1(X_{\delta}; S) - \int_0^t \sigma_u^2 \mathrm{d}u \right) \xrightarrow{Ls} MN \left( 0, \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{16}{3}c + \frac{8}{c^2} \right) \right)$$

The minimum asymptotic variance is

$$\underbrace{\frac{8\sqrt[3]{3}}{54}}_{\simeq 11.54} \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3}, \quad with \quad c = \sqrt[3]{3}$$

The Zhou estimator's asymptotic variance is thus of the form obtained by the kinked nonsubsampled realised kernels, i.e. ones which do not satisfy the k'(0) = k'(1) = 0 condition.

**Example 2** Suppose n corresponds to using prices every 1 second on the NYSE, so n = 23,400. If  $\omega^2 = 0.001$  and  $t \int_0^t \sigma_u^4 du = 1$ , which is roughly right in empirical work from 2004, then for the Zhou estimator the optimal  $S \simeq 25$  so that  $n_{\delta} \simeq 920$ . Hence the degree of subsampling is rather modest. In 2000,  $\omega^2 = 0.01$  and  $t \int_0^t \sigma_u^4 du = 1$  would be more reasonable, in which case S = 118 and  $n_{\delta} = 198$ , which corresponds to returns measured every roughly 2 minutes.

#### 3.4.2 2-lag flat-top Bartlett estimator

A natural extension of Zhou (1996) is to allow H to be larger than one but fixed.

**Lemma 4** Let  $w_0 = w_1 = 1$  and  $w_2 = 1/2$ . With  $S = c(\xi n)^{2/3}$  we have

$$n^{1/6} \left\{ \gamma_0(X_{\delta}; S) + \widetilde{\gamma}_1(X_{\delta}; S) + \frac{1}{2} \widetilde{\gamma}_2(X_{\delta}; S) - \int_0^t \sigma_u^2 \mathrm{d}u \right\} \xrightarrow{Ls} MN \left( 0, \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{20}{3} c + \frac{2}{c^2} \right) \right)$$

and the minimum variance is

$$\underbrace{10\sqrt[3]{3/5}}_{\simeq 8.43} \omega^{4/3} \left( t \int_0^t \sigma_u^4 du \right)^{2/3}, \quad with \quad c = \sqrt[3]{3/5}.$$

The constant in the asymptotic variance is here reduced from about 11.54 to 8.43. Now we proceed by adding additional realised autocovariances to Zhou's estimator, using the Bartlett weights,  $w_h = k(\frac{h-1}{H}), h = 2, ..., H$ . An interesting question is what happens as we increase H further? For moderately large H we have that  $n^{1/6} \left\{ K(X_{\delta}) - \int_0^t \sigma_u^2 du \right\}$  has an asymptotic variance of approximately  $\frac{4}{3} \left\{ 2 + (H+1) \right\} ct \int_0^t \sigma_u^4 du + \frac{8\omega^4}{c^2 H^2}$ . This suggests  $c^3 = 12\omega^4 / \left( H^3 t \int_0^t \sigma_u^4 du \right) + o(1)$ , so the asymptotic variance (using  $\frac{4}{3} 12^{1/3} + 8/12^{2/3} = 2\sqrt[3]{12}$ ) is

$$\underbrace{2\sqrt[3]{12}}_{\simeq 4.58} \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} + o(1).$$

So we achieve an additional reduction of the asymptotic variance. Not surprisingly, this is the asymptotic variance of the Bartlett realised kernel applied to a sample of size n when  $H \propto n^{2/3}$ , see Example 1. Here, by allowing H to grow we approach the situation with kinked kernels so we observe the eventual impotence of subsampling – a property we have shown holds for all kinked kernels. Hence as H gets large the optimal degree of subsampling rapidly falls and the best thing to do is simply to run a Bartlett realised kernel on the data without subsampling, i.e. take  $n_{\delta} = n$ .

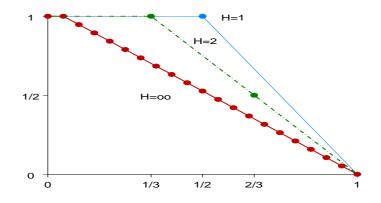


Figure 4: The implied kernels that arise from subsampling some simple kernels. H = 1 corresponds to the subsampled version of Zhou's estimator; H = 2 is that for Zhou's estimator after adding  $1/2\tilde{\gamma}_2(X_{\delta})$ ; and  $H = \infty$  (here approximated by H = 18) illustrates the implied kernel for Zhou's estimator that is enhanced by an increasing number of Bartlett-weighted realised autocovariances.

Figure 4 shows the implied kernel functions that are generated by subsampling Zhou's estimator (H = 1) and the two estimators that have been enhanced by adding Bartlett weights. The relative asymptotic efficiency for these estimators are simply given by  $k_{\bullet}^{0,0}$  of the implied kernel, where the implied kernel for H = 1 corresponds to the trapezoidal kernel by Politis and Romano (1995). From Figure 4 it is evident that  $k_{\bullet}^{0,0}$  is decreasing in H which explains that the asymptotic variance of this estimator is decreasing in H.

### 3.4.3 Relationship to two scale estimator

The two scale estimator of Zhang, Mykland, and Aït-Sahalia (2005) bias corrects  $\gamma_0(X_{\delta}; S)$  using  $\hat{\omega}^2 = \gamma_0(X_{\delta/S})/2n$ . Their results are reproved here, exploiting our previous results to make the proofs very short. We set  $S = c(\xi n)^{2/3}$ , which imposes the optimal rate for S.

**Theorem 5** With  $S = c(\xi n)^{2/3}$  we have

$$n^{1/6} \left\{ \gamma_0(X_{\delta}; S) - n_{\delta} 2\omega^2 - \int_0^t \sigma_u^2 \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left\{ 0, \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{4}{3}c + 4\frac{1+\lambda^2}{c^2} \right) \right\}, \quad (20)$$
$$n^{1/6} \left\{ \begin{array}{c} \frac{1}{S} \sum_{j=1}^{Sn_{\delta}} \left( U_{j\delta/S} - U_{(j-S)\delta/S} \right)^2 - n_{\delta} 2\omega^2 \\ \frac{1}{S} \sum_{j=1}^{Sn_{\delta}} \left( U_{j\delta/S} - U_{(j-1)\delta/S} \right)^2 - n_{\delta} 2\omega^2 \end{array} \right\} \xrightarrow{L_{\delta}} N \left\{ 0, \frac{4\omega^4}{c^2 \xi^{4/3}} \left( \begin{array}{c} 1+\lambda^2 & \lambda^2 \\ \lambda^2 & 1+\lambda^2 \end{array} \right) \right\}.$$

This allows us to understand that replacing  $\gamma_0(U_{\delta}; S) - n_{\delta} 2\omega^2$  by  $\gamma_0(U_{\delta}; S) - n_{\delta} 2\hat{\omega}^2$ , yielding a feasible estimator with a smaller variance than the infeasible estimator.

**Theorem 6** With  $S = c(\xi n)^{2/3}$  we have

$$n^{1/6} \left\{ \gamma_0(X_{\delta}; S) - 2n_{\delta}\hat{\omega}^2 - \int_0^t \sigma_u^2 \mathrm{d}u \right\} \xrightarrow{L_8} MN \left\{ 0, \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{4}{3}c + \frac{8}{c^2} \right) \right\}$$

The minimum asymptotic variance is

$$\underbrace{2\sqrt[3]{12}}_{\simeq 4.58} \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3}, \quad with \quad c = \sqrt[3]{12}.$$

Thus the two scale estimator is significantly more efficient than the Zhou estimator and is as efficient as the Bartlett realised kernel.

**Example 3** (continued from Example 2). If  $\omega^2 = 0.001$  and  $t \int_0^t \sigma_u^4 du = 1$ , then  $S \simeq 40$  and  $n_\delta \simeq 580$ . Hence the degree of subsampling is larger than that used by Zhou.

# 4 Some Empirical Recommendations

We have worked under the assumption that the noise is of the independent type defined in (6). This assumption seems reasonable for equity returns when prices are sampled at moderate high frequencies, e.g. for the liquid stocks on the New York stock exchange this assumption seems reasonable when applied to 1 minute returns (Hansen and Lunde (2006)). In this context the best approach to estimation is to use a smooth realised kernel without any subsampling. A shortcoming of this approach is that this estimator does not make use of all available observations. For example, transactions on the most liquid stocks now take place every few seconds, but for  $U \in WN$  to be reasonable we can only sample every, say, 15th observation.

In this Section we discuss how to construct subsampled realised estimators that make use of all available data. We also discuss how valid inference can be made about such estimators under realistic assumptions about the noise in tick-by-tick data.

Here we use a subsampled realised kernel, where S is chosen to be sufficiently large so that (6) is reasonable for a sample that only consists of every Sth observation. The asymptotic variance can be estimated from the coarsely sampled data, using the methods by Barndorff-Nielsen, Hansen,

Lunde, and Shephard (2006), and this leads to valid inference that is robust to both time-dependent and endogenous noise in the tick-by-tick data.

Specifically we recommend the following procedure.

- 1. Choose S sufficiently large for (6) to be a plausible assumption for a sample that only consists of every Sth observation.
- 2. Construct S distinct subsamples, each having approximately  $n_{\delta} = n/S$  observations.
- 3. For each of the S subsamples, obtain estimates of  $\omega^2$  and  $IQ = t \int_0^t \sigma_u^4 du$ , and an initial estimate of  $IV = \int_0^t \sigma_u^2 du$ . See Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) for ways to do this. Average each of these estimators to construct the subsampled estimators,  $\hat{\omega}^2 = S^{-1} \sum_{s=1}^S \hat{\omega}_s^2$  and  $\widehat{IV}_{initial} = S^{-1} \sum_{s=1}^S \widehat{IV}_{initial,s}$  and  $\widehat{IQ} = S^{-1} \sum_{s=1}^S \widehat{IQ}_s$ .
- 4. Obtain an estimate,  $\hat{H}$ , for the optimal H, by plugging the subsampled estimates into the expression for the optimal H. Use this  $\hat{H}$  to compute the S realised kernels,  $K^s(X_{\text{skip-}S})$ , using a smooth kernel and the weights  $w_0 = w_1 = 1$  and  $w_h = k\left(\frac{h-1}{\hat{H}}\right)$ , for  $h = 2, \ldots, \hat{H}$ . Form their average to obtain the actual estimator,  $\widehat{IV}_{\text{final}} = K(X_{\text{skip-}S}; S)$ .
- 5. Finally, compute the conservative estimate for avar  $\{K(X_{\text{skip-}S}; S)\}$  using the finite sample expressions, where  $w = (w_0, w_1, \dots, w_{\hat{H}})^{\intercal}$ ,

$$\widehat{\operatorname{Var}}\left\{K(X_{\operatorname{skip}-S};S)\right\} = \widehat{IQ}\left(w^{\intercal}Aw\right) \times \frac{1}{n_{\delta}} + 8\hat{\omega}^{2}\widehat{IV}_{\operatorname{final}}\left(w^{\intercal}Bw\right) + 4\hat{\omega}^{4}\left(w^{\intercal}Cw\right) \times n_{\delta}.$$
 (21)

The variance estimate in (21) is the sum of the finite sample versions of (8-10) with S = 1. So this expression completely ignores subsampling, and the expression is really an estimator of  $Var(K^s(X_{skip}))$ . The reason is that subsampling does not reduce the noise-variance by a factor of S, unless the noise is uncorrelated across subsamples. This is unrealistic when the subsamples exploit all the tick-by-tick data. However, we still have avar  $\{K(X_{skip}, S; S)\} \leq avar(K^s(X_{skip}, S))$ , even if  $U_t \perp U_s$  is violated for some  $s \neq t$ . So (21) is simply a robust estimator that is expected to yield a conservative estimate of the variance. It is interesting to have some notion of how conservative this estimator is.

Recall our result in Theorem 1 that avar  $\{K(Y_{\text{skip-}S}; S)\} = \text{avar}(K^s(Y_{\text{skip-}S}))$ , see (8). So subsampling cannot reduce the contribution to the asymptotic variance from this term, while the contributions from the two other terms (9) and (10), potentially can be driven all the way to zero.

**Example 4** With  $\rho = 1$ , the asymptotic variance of the realised TH<sub>2</sub> kernel is proportional to

$$c_1 + 2\frac{k_{\bullet,0}^{1,1}}{k_{\bullet,0}^{0,0}}c_1^{-1} + \frac{k_{\bullet,0}^{2,2}}{k_{\bullet,0}^{0,0}}c_1^{-3} = 5.75 + \frac{1.71}{0.218}\frac{2}{5.75} + \frac{41.8}{0.218}\left(5.75\right)^{-3} \simeq 9.50.$$

Subsampling this estimator with S = 10, say, reduces this factor to no less than

$$5.75 + \frac{1}{10} \frac{1.71}{0.218} \frac{2}{5.75} + \frac{1}{10} \frac{41.8}{0.218} (5.75)^{-3} \simeq 6.12,$$

see (11). So the variance reduction is less than 36% and even with  $S \to \infty$  the reduction is less than 40%. In practice, the reduction is likely to be much smaller, because the noise is not independent across subsamples. So even though (21) is a conservative estimator – it is not perversely conservative.

# 5 Simulation study

### 5.1 Simulated model and design

In this section we analyse the finite sample properties of  $K(X_{\delta}; S)$ , using both a smooth TH<sub>2</sub> kernel and a kinked Bartlett kernel. We consider the following SV model,

$$dY_t = \mu dt + \sigma_t dW_t, \quad \sigma_t = \exp\left(\beta_0 + \beta_1 \tau_t\right), \quad d\tau_t = \alpha \tau_t dt + dB_t, \quad \operatorname{corr}(dW_t, dB_t) = \rho,$$

where  $\rho$  is a leverage parameter. This model is frequently used for simulation is this context, see e.g. Huang and Tauchen (2005) and Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006).

In our simulated model, we set  $\mu = 0.03$ ,  $\beta_1 = 0.125$ ,  $\alpha = -0.025$  and  $\rho = -0.3$ . Further, we set  $\beta_0 = \beta_1^2/(2\alpha)$  in order to standardize  $E(\sigma_t^2) = 1$ . With this configuration the variance of  $\int_0^t \sigma_u^2 du$  is comparable to the empirical results found in Hansen and Lunde (2005). For the variance of market microstructure noise we set  $\omega^2 = 0.1$ .

The process is generated using an Euler scheme based on N = 23,400 intervals, where each interval is thought to correspond to one second so that the entire interval corresponds to 6.5 hours, which is the length of a typical trading day. The volatility process is restarted at its mean value  $\sigma_0 = 1$  every day by setting  $\tau_0 = 5/2$ . This keeps the noise-to-signal ratio,  $\xi = \omega^2 / \sqrt{\int_0^1 \sigma_u^4 du}$ , comparable across simulations. In our Monte Carlo designs we let the effective sample size, n, be either 1,560, 4,680, or 23,400, which correspond to sampling every 15, 5, or 1 seconds, respectively. So a sample with 4,680 observations, say, is obtained by including every fifth observation of the N = 23,401 simulated data points over the [0, t] interval.

### 5.2 Implementation of realised kernels and subsampled realised kernels

From the simulated data,  $X_0, \ldots, X_n$ , we define the "skip-S returns"  $\Delta_S X_j = X_j - X_{j-S}$ . The subsampled realised autocovariances are computed by,

$$\hat{\gamma}_{h}^{s} = \sum_{j=1}^{n_{\delta}} \Delta_{S} X_{jS+s-1} \Delta_{S} X_{(j-h)S+s-1}, \qquad s = 1, \dots, S, \qquad h = -H, \dots, 0, \dots, H,$$

where  $n_{\delta} = n/S$ . The subsampled realised kernel is defined by

$$\widehat{K(X;S)} = \frac{1}{S} \sum_{s=1}^{S} \widehat{K^s(X)}, \quad \text{where} \quad \widehat{K^s_H(X)} = \widehat{\gamma^s_0} + \sum_{h=1}^{H} k(\frac{h-1}{H}) \left(\widehat{\gamma^s_h} + \widehat{\gamma^s_{-h}}\right).$$

When S = 1 we use  $H^*_{\text{TH}_2,1} = 5.75(\xi n)^{1/2}$  for the smooth TH<sub>2</sub> kernel and  $H^*_{\text{Bartlett},1} = \sqrt[3]{12(\xi n)^2}$  for the kinked Bartlett kernel. The "noise-to-signal" parameter,  $\xi = \omega^2 / \sqrt{\int_0^1 \sigma_u^4 du}$  need not

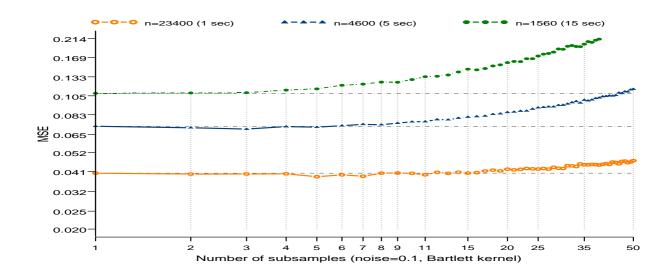


Figure 5: Mean squares errors (MSEs) for subsampled (kinked) Bartlett realised kernel using 3 different sample sizes. The MSE is fairly insensitive to S. These findings are fully consistent with Theorems 2 and 3.

be estimated in our simulations, because  $\omega^2$  is known and the integrated quarticity,  $\int_0^1 \sigma_u^4 du \simeq N \sum_{j=1}^N \sigma_{j/N}^4$ , can be computed from the simulated data. The parameter  $\rho = \int_0^1 \sigma_u^2 du/\sqrt{\int_0^1 \sigma_u^4 du}$  can be computed from the simulated volatility path. When  $S \ge 2$  the optimal H for the Bartlett kernel is simply given by  $H_{\text{Bartlett},S}^* = S^{-1} \sqrt[3]{12(\xi n)^2}$ , and the TH<sub>2</sub> kernel has  $H_{\text{TH}_2,S}^* = c_S^{1/2}(\xi n)$ , where  $c_S = S^{-1} \sqrt{7.84\rho \left(1 + \sqrt{1 + 9.33S}\right)}$ , as defined in (18).

### 5.3 Simulation Results

Figures 5 and 6 shows the Monte Carlo results with the number of subsamples, S, increasing along the horizontal axis and the MSE on the vertical axis. The lines represent different sample sizes.

Consider first the results based on the Bartlett kernel. Our theoretical results in Theorem 2 dictate that these lines should be horizontal. This result is confirmed. Still, a small increase in the MSE as S increases is observed for the smaller sample sizes. The reason is that the lag length of the implied kernel,  $H_{\text{implied}}$ , can only attain values that are divisible by S. While the Bartlett kernel without subsampling has  $H_{\text{Bartlett},1}^* = \left[\sqrt[3]{12(\xi n)^2}\right]$ , the implied Bartlett kernel has  $H_{\text{implied}} = S \times \left[S^{-1}\sqrt[3]{12(\xi n)^2}\right]$ . So as S increases the implied kernels'  $H_{\text{implied}}$  is more likely to deviate from  $H_{\text{Bartlett},1}^*$ , which causes an increase in the mean squared error. The smaller is the sample size, n, the smaller is the optimal value for H. So it is not surprising that the impact on MSE is seen earlier when n is small. In this design, the optimal lag length,  $H_{\text{Bartlett},1}^*$ , is about 67, 140, and 403, for n = 1,560, n = 4,680, and n = 23,400, respectively. Though there is some variation in the optimal H across simulations because it through  $\xi$ , depends on the simulated volatility path. The lower panels present the results for the smooth TH<sub>2</sub> kernel. Here, our theoretical results in Theorem 3 state that the MSE is increasing in S, and this phenomenon is evident for all sample

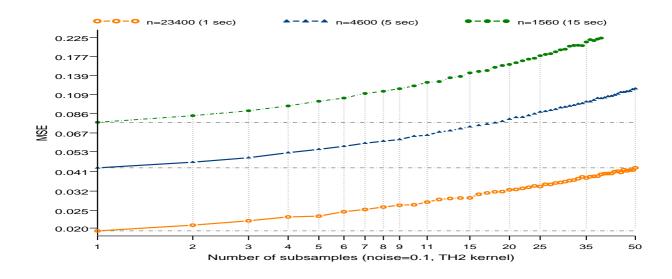


Figure 6: Mean squares errors (MSEs) for subsampled the (smooth)  $TH_2$  realised kernels using 3 different sample sizes. The  $TH_2$  kernel has MSEs that are slightly increasing in S. These findings are fully consistent Theorems 2 and 3.

sizes. The results when  $\omega^2 = 0.01$  and  $\omega^2 = 0.001$  (not reported) are similar. Here the optimal H is smaller and this causes subsampling to have a larger impact on the MSE. Naturally, the implied kernels must have  $H_{\text{implied}} \geq S$ , so that  $H_{\text{implied}} = S$  whenever  $S \geq H^*$ . This constraint is relevant for our simulations with small levels of noise because subsampling takes  $H_{\text{implied}}$  further away from its optimal value, as S increases beyond the optimal H.

# 6 Empirical study of General Electric trades

Here we compare subsampled realised kernels with other estimators. We estimate the daily increments of [Y] for the log-price of General Electric (GE) shares in 2000 and in 2004. The reason that we analyse data from both periods is that the variance of the noise was around 10 times higher in 2000 than in 2004. A more detailed analysis on 29 other major stocks is provided in a Web Appendix to this paper available from www.hha.dk/~alunde/bnhls/bnhls.htm. This appendix also describes the exact implementation of our estimators. Precise details on the cleaning we carried out on the raw data before it was analysed are described in the web appendix to Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006).

Table 3 shows Summary statistics for seven estimators. The first estimator is the realised  $TH_2$  kernel using approximate 1 minute returns. The approximate 1 minute returns are obtained by skipping a fixed number of transactions, such that the average time between observations is one minute. In 2000 we had to skip every 9.7 observations on average to construct the approximate 1 minute returns, and in 2004 we had to skip every 13.7 observations on average. The second estimator is the subsampled realised  $TH_2$  kernel. So in 2000 we have  $S \simeq 9.7$  and in 2004 we have  $S \simeq 13.7$ . The third estimator is the realised  $TH_2$  kernel that is based on tick-by-tick data (i.e. all available trades) and an H that is S times larger than that used by the first estimator.

	Mean	Std. (HAC)	$\overline{H}$	$\operatorname{Corr}$	$\operatorname{acf}(1)$	$\operatorname{acf}(2)$	$\operatorname{acf}(5)$	$\operatorname{acf}(10)$
		Sampl	e period:	2000				
Realised kernel $(TH_2)$	$H^* = ct$	$5n^{1/2})$						
$K^{\text{TH2}}(X_{\text{ap. 1 min}})$	4.747	$3.216\ (6.133)$	6.558	1.000	0.43	0.25	0.03	0.15
Subsampled realised k	ernel (T.	$H_2, H = c\xi n^{1/2})$						
$K^{{\scriptscriptstyle { m TH2}}}(X_{{ m ap. \ 1 \ min}};S)$	4.709	$3.220\ (6.170)$	6.558	0.997	0.43	0.25	0.03	0.16
Realised kernel $(TH_2,$	$H = S \cdot$	$H^*)$						
$K^{{\scriptscriptstyle{\mathrm{TH2}}}}(X_{1 \mathrm{\ tick}})$	4.702	$2.946\ (5.793)$	62.44	0.986	0.46	0.27	0.05	0.13
Subsampled realised v	ariances							
$[X_{20 \text{ minutes}}; 1200]$	4.417	$3.650\ (6.046)$		0.894	0.26	0.17	-0.01	0.17
$[X_{5 \text{ minutes}}; 300]$	4.908	3.018(5.611)		0.984	0.44	0.23	0.01	0.14
$[X_{1 \text{ minutes}}; 60]$	5.545	2.376(5.167)		0.787	0.55	0.36	0.10	0.18
AMZ (2005)								
TSRV(K, J)	4.514	3.657(6.766)		0.941	0.36	0.21	0.01	0.23
		Sampl	e period:	2004				
Realised kernel $(TH_2,$	$H^* = ct$	$(5n^{1/2})$						
$K^{{\scriptscriptstyle {\rm TH2}}}(X_{{ m ap. \ 1 \ min}})$	0.962	$0.568\ (1.195)$	5.723	1.000	0.34	0.32	0.28	0.08
Subsampled realised k	ernel (T.	$H_2, H = c\xi n^{1/2})$						
$K^{\mathrm{TH2}}(X_{\mathrm{ap.\ 1\ min}};S)$	0.954	$0.561\ (1.202)$	5.723	0.995	0.37	0.32	0.28	0.09
Realised kernel $(TH_2,$	$H = S \cdot$	$H^*)$						
$K^{\mathrm{TH2}}(X_{1 \mathrm{\ tick}})$	0.947	$0.522\ (1.130)$	78.27	0.990	0.37	0.31	0.30	0.08
Subsampled realised v	ariances							
$[X_{20 \text{ minutes}}; 1200]$	0.885	$0.516\ (1.036)$		0.933	0.27	0.27	0.27	0.08
$[X_{5 \text{ minutes}}; 300]$	0.943	0.503(1.088)		0.984	0.37	0.32	0.30	0.08
$[X_{1 \text{ minutes}}; 60]$	0.942	$0.376\ (0.921)$		0.899	0.46	0.43	0.38	0.12
AMZ (2005)								
TSRV(K, J)	0.946	$0.560\ (1.194)$		0.944	0.33	0.35	0.28	0.11

Table 3: Summary statistics for subsampled [Y] estimators.

Summary statistics for seven estimators. First the realised kernel using approximate 1 minute returns with  $H^*$  and its subsampled version, followed by the realised kernel using tick-by-tick data with  $H = S \cdot H^*$ . Then three subsampled realised variances based on 20, 5 and 1 minute returns. For instance,  $[X_{5 \text{ minutes}}; 300]$  is the average of 300 realised variances based on 5 minutes returns, obtained by shifting the time prices are recorded by 1 second. Finally, TSRV(K, J) is the two-scale estimator that is robust to deviations from i.i.d. noise. For both 2000 and 2004 we report the average of daily estimates with standard deviations. Corr is the correlation between each of estimators and the first realised kernel. Finally we report four sample autocovariances.

The following three estimators are subsampled realised variances. These are based on returns that are sampled in calendar time, where each intraday return spans 20 minutes, 5 minutes, or 1 minute, as indicated in the subscript of these estimators. To exhaust data sampled every second, the number of subsamples are S = 1200, S = 300, and S = 60, respectively. For instance, the estimator  $[X_{5 \text{ minutes}}; 300]$  is the average of 300 realised variances, where each of the realised variances are based on 5 minute intraday returns, simply changing the initial place that prices are recorded by

one second. The last estimator, TSRV(K, J), by Zhang, Mykland, and Aït-Sahalia (2005), is the two-scale estimator that is designed to be robust to deviations from i.i.d. noise. Here we use their *area adjusted* estimator, which involves a bias correction.

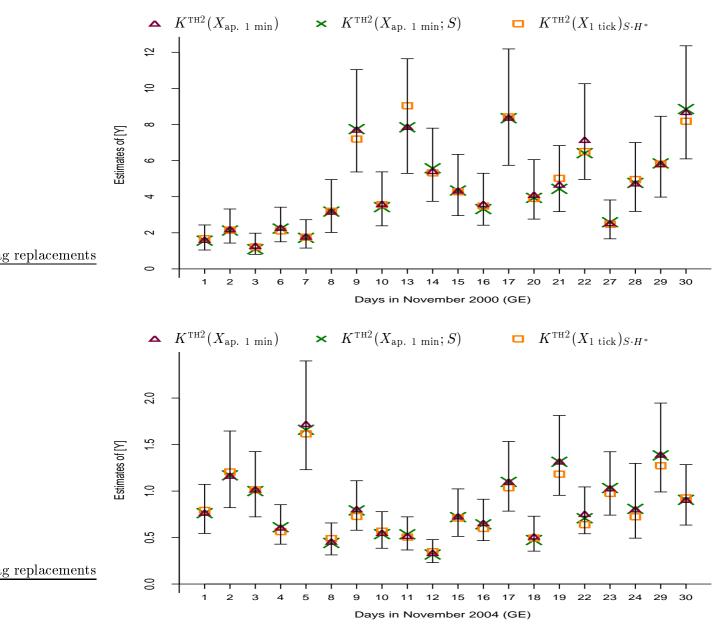


Figure 7: Three estimators for the daily increments to [Y] for General Electrics in November 2000 and 2004. Triangles are the estimates of the realised kernel using roughly 1 minute returns. Diamonds are the estimates produced by the subsampled realised kernel. Circles are the estimates of the realised kernel that uses tick-by-tick returns and an H that is S times larger than that used by the first realised kernel. The intervals are the 95% confidence intervals for  $K^{\text{TH2}}(X_{\text{ap. 1min}})$ .

From Table 3 we see that the estimators are very tightly correlated. The two realised kernels and the subsampled realised kernel are almost perfectly correlated, and all reported statistics are quite similar for these estimators. The two scale estimator is also quite similar to the realised kernels. Interestingly, amongst the subsampled realised variances, it is that based on 5 minute returns that is most similar to the realised kernels. This suggest that 20 minute returns leads to too much sampling error, whereas 1 minute returns are being influenced by the bias due to market microstructure noise.

Time series for some of these estimators are drawn in Figure 7, where we plot daily point estimates for November 2000 and November 2004. We also include the confidence intervals for  $K^{\text{TH2}}(X_{\text{ap. 1 min}})$  using the method of Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006). The three estimators are virtually almost identical. While the subsampled realised kernel may be slightly more precise than the moderately sampled realised kernel,  $K^{\text{TH2}}(X_{\text{ap. 1 min}})$ , Figure 7 does not suggest there is a big difference between these two. The realised kernel that is based on tick-by-tick data is slightly different from the other estimators, but always inside the confidence interval for  $K^{\text{TH2}}(X_{\text{ap. 1 min}})$ .

# 7 Conclusions

We have studied the properties of subsampled realised kernels. Subsampling is a very natural addition to realised kernels, for it can be viewed as averaging over realised kernels with slightly different starts of the day. We have provided a first asymptotic study of these statistics, allowing the degree of subsampling or the number of lags to go to infinity or being fixed. Included in our analysis is the asymptotic distribution of the estimator proposed by Zhou (1996).

Subsampling leads to few gains in our analysis. In fact, we found that subsampling is harmful for the best class of realised kernel estimators. The main advantage of subsampling is that it can overcome the inefficiency that results from a poor choice of kernel weights in the first place. For example, when the truncated kernel is used to design estimators, the resulting estimator has poor asymptotic properties, whereas the subsampled estimator is consistent and converges at rate  $n^{1/6}$ .

In the realistic situation where the noise is endogenous and time dependent, subsampled realised kernels do provide a simple way to make use of all the available data. We have discussed how to make valid inference about such estimators.

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# **Appendix:** Proofs

**Lemma A.1** We have  $\gamma_h(X_{\delta}; S) = \sum_{s=-S}^{S} \frac{S-|s|}{S} \gamma_{Sh+s}(X_{\frac{\delta}{S}}) + R_S^x/S.$ 

The remainder  $R_S^x/S$  is a relatively small term, due to end effects. The term is defined explicitly in the proof, and the expression shows that  $R_S^x$  can be made zero by tweaking the first S-1 and last S-1 intraday returns.

**Proof.** Define the intraday returns  $x_j = X_{\frac{\delta}{S}j} - X_{\frac{\delta}{S}j - \frac{\delta}{S}}$ , and write

$$X_{\delta(j+\frac{s-1}{S})} - X_{\delta(j-1+\frac{s-1}{S})} = X_{\frac{\delta}{S}(jS+s-1)} - X_{\frac{\delta}{S}(jS-S+s-1)} = x_{jS+s-1} + \dots + x_{jS-S+s-1}$$

So  $x_j$  are intraday returns over short intervals, each having length  $\delta/S$ . The  $\gamma_h^1(X_\delta)$  equals

$$\sum_{j=1}^{n_{\delta}} \left( X_{\delta j} - X_{\delta(j-1)} \right) \left( X_{\delta(j-h)} - X_{\delta(j-h-1)} \right)$$

$$= \sum_{j=1}^{n_{\delta}} \left( x_{(j-1)S+1} + \dots + x_{jS} \right) \left( x_{(j-h-1)S+1} + \dots + x_{(j-h)S} \right)$$

$$= \sum_{j=1}^{n} x_{j} x_{j-Sh} + \sum_{\substack{j=1\\ j \bmod S \neq 0}}^{n} x_{j} x_{j-Sh+1} + \sum_{\substack{j=1\\ j \bmod S \notin \{0,S-1\}}}^{n} x_{j} x_{j-Sh+2} + \dots + \sum_{\substack{j=1\\ j \bmod S=1}}^{n} x_{j} x_{j-Sh+S-1}$$

$$+ \sum_{\substack{j=1\\ j \bmod S \neq 1}}^{n} x_{j} x_{j-Sh-1} + \sum_{\substack{j=1\\ j \bmod S \notin \{1,2\}}}^{n} x_{j} x_{j-Sh-2} + \dots + \sum_{\substack{j=1\\ j \bmod S=0}}^{n} x_{j} x_{j-Sh-S+1}.$$

Similarly for s > 1 we have

$$\begin{split} &\sum_{j=1}^{n_{\delta}} \left( X_{\delta j + \frac{s-1}{S}} - X_{\delta (j-1) + \frac{s-1}{S}} \right) \left( X_{\delta (j-h) + \frac{s-1}{S}} - X_{\delta (j-h-1) + \frac{s-1}{S}} \right) = \sum_{j=s}^{n+s-1} x_j x_{j-Sh} \\ &+ \sum_{\substack{j=s \\ j \bmod S \neq s-1}}^{n+s-1} x_j x_{j-Sh+1} + \sum_{\substack{j=s \\ j \bmod S \notin \{s-1,s-2\}}}^{n+s-1} x_j x_{j-Sh+2} + \dots + \sum_{\substack{j=s \\ j \bmod S = s}}^{n+s-1} x_j x_{j-Sh+S-1} \\ &+ \sum_{\substack{j=s \\ j \bmod S \neq s}}^{n+s-1} x_j x_{j-Sh-1} + \sum_{\substack{j=s \\ j \bmod S \notin \{s,1\}}}^{n+s-1} x_j x_{j-Sh-2} + \dots + \sum_{\substack{j=s \\ j \bmod S = s-1}}^{n+s-1} x_j x_{j-Sh-S+1} . \end{split}$$

By adding up the terms,  $\gamma_h(X_{\delta}; S) = \sum_{s=-S+1}^{S-1} \frac{S-s}{S} \gamma_{Sh+s}(X_{\frac{\delta}{S}}) + R_S^x/S$ , where

$$R_{S}^{x} = -\sum_{S=2}^{S} \qquad \left(\sum_{j=1}^{s-1} x_{j} x_{j-Sh} + \sum_{j=1}^{s-2} x_{j} x_{j-Sh+1} + \dots + \sum_{j=1}^{1} x_{j} x_{j-Sh+S-2} + \dots + \sum_{j=1}^{s-2} x_{j} x_{j-Sh+1} + \dots + \sum_{j=1}^{s-2} x_{j-Sh+1} + \dots + \sum_{j=1}^$$

$$+\sum_{j=1}^{s-1} x_j x_{j-Sh-1} + \sum_{j=2}^{s-1} x_j x_{j-Sh-2} + \dots + \sum_{j=s-1}^{s-1} x_j x_{j-Sh-S+1} \right)$$
  
+ 
$$\sum_{S=2}^{S} \left( \sum_{j=n+1}^{n+s-1} x_j x_{j-Sh} + \sum_{j=n+1}^{n+s-2} x_j x_{j-Sh+1} + \dots + \sum_{j=n+1}^{n+1} x_j x_{j-Sh+S-2} + \sum_{j=n+1}^{n+s-1} x_j x_{j-Sh-1} + \sum_{j=n+2}^{n+s-1} x_j x_{j-Sh-2} + \dots + \sum_{j=n+s-1}^{n+s-1} x_j x_{j-Sh-h+1} \right).$$

The term,  $R_S^x$ , is due to end effects and involves much fewer cross products,  $x_i x_j$ , than does  $\sum_{s=1}^{S} \gamma_h^s(X_\delta)$ . So that  $R_S^x/S$  is typically negligible. In fact,  $R_S^x$  can be made zero by assuming  $x_1 = \cdots = x_{S-1} = x_{n+1} = \cdots = x_{n+S-1} = 0$ .  $\Box$ 

In the non-subsampling case Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) derived the following helpful results.

**Theorem A.1** We study properties as  $\delta \downarrow 0$ . Suppose that  $Y \in \mathcal{BSM}$  and (4) holds, then

$$n_{\delta}^{1/2} \begin{pmatrix} \gamma_0(Y_{\delta}) - \int_0^t \sigma_u^2 \mathrm{d}u \\ \tilde{\gamma}_1(Y_{\delta}) \\ \vdots \\ \tilde{\gamma}_H(Y_{\delta}) \end{pmatrix} \xrightarrow{Ls} MN \left( 0, A_1 \times t \int_0^t \sigma_u^4 \mathrm{d}u \right), \quad A_1 = \begin{pmatrix} 2 & 0 & \cdots & 0 \\ 0 & 4 & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & 4 \end{pmatrix}.$$
(A.1)

If, in addition, for  $n_{\delta} \geq H$ ,  $U \in \mathcal{WN}$  and  $Y \perp U$  then  $\widetilde{\gamma}(Y_{\delta}, U_{\delta}) \xrightarrow{L_{\delta}} MN(0, 2\omega^{2}[Y]B)$ , where

$$B = \begin{pmatrix} B_{11} & B_{12} \\ B_{21} & B_{22} \end{pmatrix}, \ B_{22} = \begin{pmatrix} 2 & -1 & 0 & \cdots \\ -1 & 2 & \ddots & \ddots \\ \ddots & \ddots & \ddots & -1 \\ \cdots & 0 & -1 & 2 \end{pmatrix}, \ B_{11} = \begin{pmatrix} 1 & \bullet \\ -1 & 2 \end{pmatrix}, \ B_{21} = \begin{pmatrix} 0 & -1 \\ 0 & 0 \\ \vdots & \vdots \\ 0 & 0 \end{pmatrix},$$

$$E \{ \widetilde{\gamma}(U_{\delta}) \} = 2\omega^{2} n_{\delta} (1, -1, 0, 0, ..., 0)^{\mathsf{T}}, \quad and \quad \operatorname{Cov} \{ \widetilde{\gamma}(U_{\delta}) \} = 4\omega^{4} n_{\delta} C + O(1),$$

$$C = \begin{pmatrix} C_{11} & C_{12} \\ C_{21} & C_{22} \end{pmatrix}, \quad C_{11} = \begin{pmatrix} 1 + \lambda^{2} & -2 - \lambda^{2} \\ -2 - \lambda^{2} & 5 + \lambda^{2} \end{pmatrix},$$

$$C_{21} = \begin{pmatrix} 1 & -4 \\ 0 & 1 \\ 0 & 0 \\ \vdots & \vdots \\ 0 & 0 \end{pmatrix}, \quad C_{22} = \begin{pmatrix} 6 & \bullet & \bullet & \bullet \\ -4 & 6 & \bullet & \bullet \\ 1 & -4 & 6 & \bullet \\ 0 & 1 & -4 & 6 & \bullet \\ \vdots & \ddots & \ddots & \ddots & \ddots \end{pmatrix}.$$

**Theorem A.2** Suppose that  $Y \in \mathcal{BSM}$  and (4) holds, then as  $\delta \downarrow 0$ 

$$n_{\delta}^{1/2} \begin{pmatrix} \gamma_0(Y_{\delta}; S) - \int_0^t \sigma_u^2 \mathrm{d}u, \\ \widetilde{\gamma}_1(Y_{\delta}; S) \\ \vdots \\ \widetilde{\gamma}_H(Y_{\delta}; S) \end{pmatrix} \xrightarrow{Ls} MN\left(0, A_S \times t \int_0^t \sigma_u^4 \mathrm{d}u\right),$$

$$A_{S} = \frac{2}{3} \begin{pmatrix} 2+S^{-2} & \bullet & 0 & \cdots \\ 1-S^{-2} & 4+2S^{-2} & \bullet & \ddots \\ 0 & 1-S^{-2} & 4+2S^{-2} & \ddots \\ \vdots & \ddots & \ddots & \ddots \end{pmatrix} \rightarrow \frac{2}{3} \begin{pmatrix} 2 & 1 & 0 & \cdots \\ 1 & 4 & 1 & \ddots \\ 0 & 1 & 4 & \ddots \\ \vdots & \ddots & \ddots & \ddots \end{pmatrix} = A_{\infty}, \quad (A.2)$$

and as  $\delta \downarrow 0$  and  $S \rightarrow \infty$ 

$$n_{\delta}^{1/2} \begin{pmatrix} \gamma_0(Y_{\delta}; S) - \int_0^t \sigma_u^2 \mathrm{d}u, \\ \widetilde{\gamma}_1(Y_{\delta}; S) \\ \vdots \\ \widetilde{\gamma}_H(Y_{\delta}; S) \end{pmatrix} \xrightarrow{Ls} MN \left( 0, A_{\infty} \times t \int_0^t \sigma_u^4 \mathrm{d}u \right)$$

**Proof of Theorem A.2.** By Lemma A.1 we have  $\tilde{\gamma}_h(Y_{\delta}; S) \simeq \sum_{s=-S}^{S} \frac{S-|s|}{S} \tilde{\gamma}_{Sh+s}(Y_{\delta})$ , and the asymptotic properties of  $\gamma_h(Y_{\delta})$ ,  $h = -SH, \ldots, SH$ , using the small time gaps,  $\delta/S$ , follows straightforwardly from (A.1). Write

$$V_{0,S} = \frac{1}{S} \left( 1 + 2\sum_{s=1}^{S} \left( \frac{S-s}{S} \right)^2 \right) = \frac{2}{3} \left( 1 + \frac{S^{-2}}{2} \right) \to \frac{2}{3} \quad V_{1,S} = \frac{1}{S} \left( 0 + \sum_{s=1}^{S} \frac{s}{S} \frac{S-s}{S} \right) = \frac{1}{6} \left( 1 - S^{-2} \right) \to \frac{1}{6}$$

then for  $h \ge 1$  we have

$$\operatorname{Var}\left\{\tilde{\gamma}_{h}(Y_{\delta};S)\right\} = \operatorname{Var}\left\{\sum_{s=-S}^{S} \frac{S-|s|}{S}\tilde{\gamma}_{Sh+s}(Y_{\delta_{S}})\right\} \to \frac{4V_{0,S}}{n_{\delta}}t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u_{S}^{4}$$

and similarly for h = 0 we find  $\operatorname{Var} \{ \tilde{\gamma}_0(Y_{\delta}; S) \} = \frac{2V_{0,S}}{n_{\delta}} t \int_0^t \sigma_u^4 \mathrm{d}u$ . For  $h \ge 0$  we find

$$\begin{aligned} \operatorname{Cov}\left\{\tilde{\gamma}_{h}(Y_{\delta};S),\tilde{\gamma}_{h+1}(Y_{\delta};S)\right\} &= \operatorname{Cov}\left\{\sum_{s=-S}^{S} \frac{S-|s|}{S}\tilde{\gamma}_{Sh+s}(Y_{\delta_{S}}), \sum_{s=-S}^{S} \frac{S-|s|}{S}\tilde{\gamma}_{Sh+S+s}(Y_{\delta_{S}})\right\} \\ &= \operatorname{Var}\left\{\sum_{s=1}^{S} \frac{S-s}{S}\tilde{\gamma}_{Sh+s}(Y_{\delta_{S}})\right\} = \sum_{s=1}^{S} \frac{S-s}{S}\frac{s}{S} \times \frac{1}{n} 4t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u \\ &= 4V_{1,S} \times \frac{1}{n_{\delta}} t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u.\end{aligned}$$

Covariances between  $\tilde{\gamma}_h(Y_{\delta}; S)$  and  $\tilde{\gamma}_i(Y_{\delta}; S)$  are zero for  $|h - i| \ge 2$ , as they do not "share" any of the realised autocovariances  $\tilde{\gamma}_{Sh+s}(Y_{\delta_S})$ .  $\Box$ 

**Proof of Theorem 1.** For the subsampled realised kernel on  $Y_{\delta}$  we have

$$S(V_{0,S} + 2V_{1,S}) = 1 + 2\sum_{s=1}^{S} \left(\frac{S-s}{S}\right)^2 + 2\sum_{s=1}^{S} \frac{s}{S} \frac{S-s}{S} + \frac{S(S-1)}{S} = S,$$

so that  $V_{0,S} + 2V_{1,S} = 1$ , where  $V_{0,S}$  and  $V_{1,S}$  are defined in the proof of Theorem A.2. From the structure of  $A_S$  we have

$$\frac{1}{H} \sum_{i,j=0}^{H} k(\frac{i}{H}) k(\frac{j}{H}) [A_S]_{i,j} = \frac{4V_{0,S}}{H} \sum_{h=0}^{H} k(\frac{h}{H})^2 + \frac{8V_{1,S}}{H} \sum_{h=1}^{H} k(\frac{h}{H}) k(\frac{h-1}{H}) + O(\frac{1}{H})$$

$$= \frac{4(V_{0,S} + 2V_{1,S})}{H} \sum_{h=0}^{H} k(\frac{h}{H})^2 - \frac{8V_{1,S}}{H^2} \sum_{h=1}^{H} k(\frac{h}{H}) \frac{k(\frac{h}{H}) - k(\frac{h-1}{H})}{1/H} + O(\frac{1}{H}) = 4 \int_0^1 k(u)^2 dx + O(\frac{1}{H}).$$

From Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) we have

$$\widetilde{\gamma}(Y_{\delta}, U_{\delta}; S) \xrightarrow{L_{\delta}} MN\left(0, \frac{2\omega^2}{S}[Y]B\right),$$
(A.3)

$$E\{\tilde{\gamma}(U_{\delta};S)\} = 2\omega^2 n_{\delta} (1, -1, 0, 0, ..., 0)^{\mathsf{T}}.$$
(A.4)

Furthermore, with  $U_j^s = U_{(j+\frac{s-1}{S})\delta}$  we have  $\tilde{\gamma}_h^s(U_{\delta}) = -V_{h+1,n}^s + 2V_{h,n}^s - V_{h-1,n}^s + R_{h+1,n}^s - R_{h-1,n}^s$ , where  $V_{h,n}^s = \sum_{j=1}^n U_j^s(U_{j-h}^s + U_{j+h}^s)$  and  $R_{h,n}^s = \frac{1}{2}(U_n^s U_{n+h}^s + U_0^s U_{-h}^s - U_n^s U_{n-h}^s - U_0 U_h^s)$ . So with  $w_0 = w_{-1} = 1$ ,  $w_h = k(\frac{h-1}{H})$ ,  $h = 1, \ldots, H + 1$  we have from Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006) that

$$K^{s}(U_{\delta}) = -\sum_{h=0}^{H} \left( w_{h+1} - 2w_{h} + w_{h-1} \right) V_{h,n}^{s} - \sum_{h=1}^{H} \left( w_{h+1} - w_{h-1} \right) R_{h,n}^{s},$$

where  $\left[\frac{n}{H^3}k_{\bullet}^{2,2} + \frac{n}{H^2}\{k'(0) + k'(1)\}\right]^{-1/2} \sum_{h=0}^{H} (w_{h+1} - 2w_h + w_{h-1}) V_{h,n}^s \xrightarrow{d} N(0, 4\omega^4)$  and  $\operatorname{avar}(H^{-1/2}\sum_{h=1}^{H} (w_{h+1} - w_{h-1}) R_{h,n}^s) = 4\omega^4 k_{\bullet}^{1,1}$ . Since the noise is independent across subsamples, the results for  $K(X_{\delta}, S) + \Delta_{H,n}^S = -\sum_{h=1}^{H} (w_{h+1} - 2w_h + w_{h-1}) S^{-1} \sum_{s=1}^{S} V_{h,n}^s$  and  $\Delta_{H,n}^S = \sum_{h=1}^{H} (w_{h+1} - w_{h-1}) S^{-1} \sum_{s=1}^{S} R_{h,n}^s$  follow.  $\Box$ 

**Proof of Lemma 1.** From (12) we have

$$\frac{\operatorname{avar}\{K(X_{\delta}) - \tilde{K}(X_{\delta})\}}{\operatorname{avar}\{\tilde{K}(X_{\delta})\}} = \frac{4\omega^{4}k_{\bullet}^{1,1}/(HS)}{4t\int_{0}^{t}\sigma_{u}^{4}\mathrm{d}u\left[\frac{HS}{n}k_{\bullet}^{0,0} + \frac{2\xi\rho k_{\bullet}^{1,1}}{HS} + n\xi^{2}\left\{\frac{k'(0)^{2} + k'(1)^{2}}{(HS)^{2}} + S\frac{k_{\bullet}^{2,2}}{(HS)^{3}}\right\}\right]}{\frac{\xi}{\frac{(HS)^{2}}{k_{\bullet}^{1,1}} + 2\rho + \frac{n}{HS}\xi\frac{k'(0)^{2} + k'(1)^{2}}{k_{\bullet}^{1,1}} + S\frac{n}{(HS)^{2}}\xi\frac{k_{\bullet}^{2,2}}{k_{\bullet}^{1,1}}},$$

which can be seen to vanish when  $k'(0)^2 + k'(1)^2 \neq 0$  or  $S \to \infty$ . We need  $HS \propto n^{1/2}$  for the ratio not to vanish when  $k'(0)^2 + k'(1)^2 = 0$ . With  $HS = c\xi \sqrt{n}$  we find

$$\frac{\operatorname{avar}\{K(X_{\delta}) - \tilde{K}(X_{\delta})\}}{\operatorname{avar}\{\tilde{K}(X_{\delta})\}} = \frac{\xi}{c^{2}\xi \frac{k_{\bullet}^{0,0}}{k_{\bullet}^{1,1}} + 2\rho + \frac{S}{\xi c^{2}} \frac{k_{\bullet}^{2,2}}{k_{\bullet}^{1,1}}} \le \frac{\xi}{2\left(1 + \sqrt{\frac{k_{\bullet}^{2,2}k_{\bullet}^{0,0}}{(k_{\bullet}^{1,1})^{2}}}\right)},$$

where we used that  $\rho, S \ge 1$  and that  $x = \sqrt{b/a}$  minimizes f(x) = ax + b/x, a, b > 0. **Proof of Theorem 2.** (i) The mixed Gaussian result follows from Theorem 1. (ii) The best value for c is found by solving the first order condition  $k_{\bullet}^{0,0} - 2c^{-3} \{k'(0)^2 + k'(1)^2\} = 0$ , and substituting this c into (13) yields  $\omega^{4/3} \left(t \int_0^t \sigma_u^4 du\right)^{2/3}$  times

$$4c\left\{k_{\bullet}^{0,0} + \frac{k'(0)^2 + k'(1)^2}{c^3}\right\} = 4c\left(k_{\bullet}^{0,0} + \frac{1}{2}k_{\bullet}^{0,0}\right) = 4ck_{\bullet}^{0,0}\left(1 + 1/2\right) = 6ck_{\bullet}^{0,0}.$$

Finally  $ck_{\bullet}^{0,0} = \left\{ 2 \left( k'(0)^2 + k'(1)^2 \right) / k_{\bullet}^{0,0} \right\}^{1/3} k_{\bullet}^{0,0} = \left\{ 2 \left( k_{\bullet}^{0,0} \right)^2 \left( k'(0)^2 + k'(1)^2 \right) \right\}^{1/3} . \square$ **Proof of Theorem 3.** (*i.a*) The mixed Gaussian result is straight forward using Theorem 1.

(*i.b*) Substituting  $HS = \xi^{1/2} c n^{1/2 + \alpha/4}$  and  $S = a n^{\alpha}$  into (14) yields  $4\omega \left( t \int_0^t \sigma_u^4 du \right)^{3/4}$  times

$$\frac{cn^{1/2+\alpha/4}}{n}k_{\bullet}^{0,0} + \frac{2\rho k_{\bullet}^{1,1}}{cn^{1/2+\alpha/4}} + nn^{\alpha}\frac{k_{\bullet}^{2,2}}{(cn^{1/2+\alpha/4})^3} = ck_{\bullet}^{0,0}n^{-1/2+\alpha/4} + c^{-3}k_{\bullet}^{2,2}n^{-1/2+\alpha/4},$$

because the second term is of lower order that the 1st and 3rd term when  $\alpha > 0$ .

(ii) Minimizing (14) with respect to x = HS has the first order condition,

$$n^{-1}k_{\bullet}^{0,0} - 2\xi\rho k_{\bullet}^{1,1}(HS)^{-2} - 3\xi^2 nSk_{\bullet}^{2,2}(HS)^{-4} = 0.$$

The unique positive solution is given by  $HS = c_S(\xi n)^{1/2}$ , where

$$c_{S} = \sqrt{\frac{\rho k_{\bullet}^{1,1}}{k_{\bullet}^{0,0}} \left(1 + \sqrt{1 + 3S \frac{k_{\bullet}^{0,0} k_{\bullet}^{2,2}}{(\rho k_{\bullet}^{1,1})^{2}}}\right)} = \sqrt{\frac{\rho k_{\bullet}^{1,1}}{k_{\bullet}^{0,0}} + \sqrt{\frac{(\rho k_{\bullet}^{1,1})^{2} + 3S k_{\bullet}^{0,0} k_{\bullet}^{2,2}}{(k_{\bullet}^{0,0})^{2}}}.$$

Now define  $x = k_{\bullet}^{0,0}/(\rho k_{\bullet}^{1,1}), y = \rho k_{\bullet}^{1,1}/(Sk_{\bullet}^{2,2})$ , and  $z = \sqrt{1+3x/y}$ . Then  $c_S = \sqrt{(1+z)/x}$ and  $x/y = (z^2 - 1)/3 = (1+z)(z-1)/3$ . So the minimum asymptotic variance is given by  $4\omega \left(t \int_0^t \sigma_u^4 du\right)^{3/4} k_{\bullet}^{0,0} (c_S + \frac{2}{c_S x} + \frac{1}{c_S^3 xy})$ , which is proportional to

$$c_S + \frac{2}{c_S x} + \frac{1}{c_S^3 x y} = \sqrt{\frac{1+z}{x}} + 2\sqrt{\frac{1}{x(1+z)}} + \frac{\sqrt{x}}{(1+z)\sqrt{1+z}c^3 y} = \frac{1}{\sqrt{x}} \frac{4}{3} \left(\frac{1}{\sqrt{1+z}} + \sqrt{1+z}\right).$$

Now substitute  $z = \sqrt{1 + 3Sk_{\bullet}^{0,0}k_{\bullet}^{2,2}/(\rho k_{\bullet}^{1,1})^2}$  and  $x^{-1/2} = \sqrt{\rho k_{\bullet}^{1,1}/k_{\bullet}^{0,0}}$  and (16) follows.  $\Box$ 

**Lemma A.2** Let g(S) be as defined in Theorem 3. Then g'(S) > 0 for all S > 0.

**Proof.** Consider the function  $f(x) = \frac{1}{\sqrt{1+\sqrt{1+ax}}} + \sqrt{1+\sqrt{1+ax}}$ , for a > 0. The first derivative  $f'(x) = \frac{a}{4} \left(1 + \sqrt{ax+1}\right)^{-3/2}$ , is positive for all x > 0.  $\Box$ 

**Proof of Corollary 1.** From Lemma A.2 it follows that g'(S) > 0 for all S > 0, if we set x = S and  $a = 3k_{\bullet}^{0,0}k_{\bullet}^{2,2}/(\rho k_{\bullet}^{1,1})^2$ . So any increment in S will increase the asymptotic variance.  $\Box$ 

**Proof of Corollary 2.** By substitution for the first  $\rho$  in g(S) we find that (16) is proportional to

$$\omega \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{1/2} \left( \int_0^t \sigma_u^2 \mathrm{d}u \right)^{1/2} \left\{ \frac{1}{\sqrt{1 + \sqrt{1 + 3Sk_{\bullet}^{0,0}k_{\bullet}^{2,2}/(\rho k_{\bullet}^{1,1})^2}} + \sqrt{1 + \sqrt{1 + 3Sk_{\bullet}^{0,0}k_{\bullet}^{2,2}/(\rho k_{\bullet}^{1,1})^2} \right\}.$$

From Hansen and Lunde (2006, p. 135) it follows that business time sampling minimizes  $t \int_0^t \sigma_u^4 du$ and by Lemma A.2 we have that also the second term is minimized for the largest possible value of  $\rho$ , (set  $x = 1/\rho^2$ ). Since  $\rho \leq 1$  the solution is  $\rho = 1$ .  $\Box$ 

**Proof of Lemma 2.** From the proof of 1 we have

$$K_w(U_{\delta}) = K_w(U_{\delta}; 1) = -\sum_{h=0}^{H} (w_{h+1} - 2w_h + w_{h-1}) V_{h,n}^1 - \sum_{h=1}^{H} (w_{h+1} - w_{h-1}) R_{h,n}^1$$

where  $\operatorname{Var}(V_{h,n}^1) = (4n-2h)\omega^4$ .  $V_{h,n}^1$  is entirely made up of  $U_j^1 U_{j-h}^1$  terms so that  $\operatorname{Cov}(V_{h,n}^1, V_{k,n}^1) = 0$ , for  $h \neq k$ . Hence  $\operatorname{Var}\left\{\tilde{K}_w(U_\delta)\right\} \geq 4\omega^4 (n-\frac{H}{2})\sum_{h=0}^H (w_{h+1}-2w_h+w_{h-1})^2$ , and the results follows since H = o(n).  $\Box$ 

**Proof of Theorem 4.** The asymptotic distribution of  $\gamma_0(X_{\delta}; S) + \tilde{\gamma}_1(X_{\delta}; S) - \int_0^t \sigma_u^2 du$  is mixed Gaussian with variance of approximately, for moderate  $n_{\delta}$  and S,

$$n_{\delta}^{-1} \frac{16}{3} t \int_{0}^{t} \sigma_{u}^{4} \mathrm{d}u + \frac{8\omega^{4} n_{\delta}}{S}.$$
 (A.5)

The first term appears from (A.2), the second from Theorem A.2 of Barndorff-Nielsen, Hansen, Lunde, and Shephard (2006).  $\Box$ 

**Proof of Lemma 3.** With  $S = c(\xi n)^{2/3}$  we have  $n_{\delta}^{-1} = S/n = c\xi^{2/3}n^{-1/3}$  and  $\frac{n_{\delta}}{S} = n/S^2 = c^{-2}\xi^{-4/3}n^{1/3}$ , so that (A.5) in the proof of Theorem 4 becomes  $n^{1/3}$  times

$$\frac{16}{3}c\xi^{2/3}t\int_0^t \sigma_u^4 \mathrm{d}u + 8\omega^4 c^{-2}\xi^{-4/3} = \omega^{4/3} \left(t\int_0^t \sigma_u^4 \mathrm{d}u\right)^{2/3} \left(\frac{16}{3}c + 8c^{-2}\right).$$

So  $n^{1/6} \left\{ \gamma_0(X_{\delta}; S) + \tilde{\gamma}_1(X_{\delta}; S) - \int_0^t \sigma_u^2 du \right\}$  converges to a mixed Gaussian distribution with this variance. We can now minimise this asymptotic variance by selecting  $c^3 = 3$ . At this value the asymptotic variance is

$$\omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left\{ \frac{16}{3} \left( 3 \right)^{1/3} + 8 \left( 3 \right)^{-2/3} \right\} \simeq 11.53 \omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3}.$$

**Proof of Lemma 4.** From Theorems A.2 and 1 we obtain the following upper-left  $3 \times 3$  submatrices of  $A_{\infty}$  and C,

$$[A_{\infty,3\times3}] = \frac{2}{3} \begin{pmatrix} 2 & \bullet & \bullet \\ 1 & 4 & \bullet \\ 0 & 1 & 4 \end{pmatrix}, \qquad [C_{3\times3}] = \begin{pmatrix} \lambda^2 + 1 & \bullet & \bullet \\ -\lambda^2 - 2 & \lambda^2 + 5 & \bullet \\ 1 & -4 & 6 \end{pmatrix}.$$

With  $w = (1, 1, \frac{1}{2})^{\mathsf{T}}$  we have  $w^{\mathsf{T}}[A_{\infty,3\times3}]w = \frac{20}{3}$  and  $w^{\mathsf{T}}[C_{3\times3}]w = \frac{1}{2}$ . The result now follows, as the asymptotic variance is

$$n^{1/3} \left( \frac{S}{n} t \int_0^1 \sigma_u^4 \mathrm{d}u \frac{20}{3} + 4\omega^4 \frac{n}{S^2} \frac{1}{2} \right) = \omega^{4/3} \left( t \int_0^1 \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{20}{3} c + 2c^{-2} \right)$$

### Proof of Theorem 5. We have

$$\gamma_0(X_{\delta};S) - 2n_{\delta}\omega^2 - \int_0^t \sigma_u^2 \mathrm{d}u = \underbrace{\gamma_0(Y_{\delta};S) - \int_0^t \sigma_u^2 \mathrm{d}u}_{n_{\delta}^{-1}\frac{4}{3}t \int_0^t \sigma_u^4 \mathrm{d}u} + \underbrace{2\gamma_0(U_{\delta},Y_{\delta};S)}_{S^{-1}8\omega^2 \int_0^t \sigma_u^2 \mathrm{d}u} + \underbrace{\gamma_0(U_{\delta};S) - 2n_{\delta}\omega^2}_{4\omega^4 \frac{n_{\delta}}{S}(1+\lambda^2)},$$

which has mean zero and a variance that is the sum of the three terms given below the brackets. The three terms are given from (A.2), (A.3), and Theorem A.1, respectively. For large  $S = c(\xi n)^{2/3}$ (implying large  $n_{\delta} = n/S = c^{-1}\xi^{-2/3}n^{1/3}$ ) we have

$$n^{1/6} \left\{ \gamma_0(X_{\delta}; S) - 2n_{\delta}\omega^2 - \int_0^t \sigma_u^2 \mathrm{d}u \right\} \xrightarrow{L_{\delta}} MN \left\{ 0, 4\omega^{4/3} \left( t \int_0^t \sigma_u^4 \mathrm{d}u \right)^{2/3} \left( \frac{c}{3} + \frac{1+\lambda^2}{c^2} \right) \right\}.$$

By the approximations

$$\frac{1}{S} \sum_{j=1}^{n} \left( U_{j\delta/S} - U_{(j-S)\delta/S} \right)^2 \simeq \frac{2}{S} \left( \sum_{j=1}^{n} U_{j\delta/S}^2 + \sum_{j=1}^{n} U_{j\delta/S} U_{(j-S)\delta/S} \right)$$
$$\frac{1}{S} \sum_{j=1}^{n} \left( U_{j\delta/S} - U_{(j-1)\delta/S} \right)^2 \simeq \frac{2}{S} \left( \sum_{j=1}^{n} U_{j\delta/S}^2 + \sum_{j=1}^{n} U_{j\delta/S} U_{(j-1)\delta/S} \right),$$

and using  $\frac{2}{S} = \frac{2n^{1/2}}{c\xi^{2/3}n^{2/3}} \times n^{-1/2} = n^{-1/6}\sqrt{\frac{4}{c^2\xi^{4/3}}} \times n^{-1/2}$  we see that

$$n^{-1/6} \left( \begin{array}{c} n^{-1/2} \sum_{j=1}^{n} \left( U_{j\delta/S} - U_{(j-S)\delta/S} \right)^2 - 2n_{\delta}\omega^2 \\ n^{-1/2} \sum_{j=1}^{n} \left( U_{j\delta/S} - U_{(j-1)\delta/S} \right)^2 - 2n_{\delta}\omega^2 \end{array} \right) \xrightarrow{L} N \left\{ 0, \frac{4\omega^4}{c^2 \xi^{4/3}} \left( \begin{array}{c} 1 + \lambda^2 & \lambda^2 \\ \lambda^2 & 1 + \lambda^2 \end{array} \right) \right\}.$$

**Proof of Theorem 6.** Follows from Theorem 5, and  $n^{-1/2}\gamma_0(X_{\delta/S}) = n^{-1/2}\gamma_0(U_{\delta/S}) + o_p(1)$  and  $\omega^4/\xi^{4/3} = \omega^{4/3} \left(t \int_0^t \sigma_u^4 du\right)^{2/3}$ .  $\Box$