

# Minimax and Adaptive Prediction for Functional Linear Regression

T. Tony CAI and Ming YUAN

This article considers minimax and adaptive prediction with functional predictors in the framework of functional linear model and reproducing kernel Hilbert space. Minimax rate of convergence for the excess prediction risk is established. It is shown that the optimal rate is determined jointly by the reproducing kernel and the covariance kernel. In particular, the alignment of these two kernels can significantly affect the difficulty of the prediction problem. In contrast, the existing literature has so far focused only on the setting where the two kernels are nearly perfectly aligned. This motivates us to propose an easily implementable data-driven roughness regularization predictor that is shown to attain the optimal rate of convergence adaptively without the need of knowing the covariance kernel. Simulation studies are carried out to illustrate the merits of the adaptive predictor and to demonstrate the theoretical results.

KEY WORDS: Functional linear model; Minimax rate of convergence; Principal components analysis; Reproducing kernel Hilbert space; Spectral decomposition.

## 1. INTRODUCTION

Prediction is an important problem in functional data analysis with a wide range of applications including chemometrics, econometrics, and biomedical studies. See, for example, Ramsay and Silverman (2002, 2005), Ferraty and Vieu (2006), and Ramsay, Hooker, and Graves (2009). Consider the functional linear model

$$Y = \alpha_0 + \int_{\mathcal{T}} X(t)\beta_0(t)dt + \epsilon, \quad (1)$$

where  $Y$  is a scalar response,  $X : \mathcal{T} \rightarrow \mathbb{R}$  is a square integrable functional predictor defined over a compact domain  $\mathcal{T} \subset \mathbb{R}$ ,  $\alpha_0$  is the intercept,  $\beta_0 : \mathcal{T} \rightarrow \mathbb{R}$  is the slope function, and  $\epsilon$  is the random noise with mean 0 and finite variance  $\sigma^2$ . In this article, we focus on the random design where  $X$  is a path of a square integrable stochastic process defined over  $\mathcal{T}$  and is independent of  $\epsilon$ . The goal of prediction is to recover the functional  $\eta_0$ :

$$\eta_0(X) = \alpha_0 + \int_{\mathcal{T}} X(t)\beta_0(t)dt,$$

based on a training sample  $\{(X_i, Y_i) : i = 1, \dots, n\}$  consisting of  $n$  independent copies of  $(X, Y)$ . Let  $\hat{\eta}_n$  be a prediction rule constructed from the training data. Then, its accuracy can be naturally measured by the excess risk:

$$\begin{aligned} \mathcal{E}(\hat{\eta}_n) &:= \mathbb{E}^* [Y^* - \hat{\eta}_n(X^*)]^2 - \mathbb{E}^* [Y^* - \eta_0(X^*)]^2 \\ &= \mathbb{E}^* [\hat{\eta}_n(X^*) - \eta_0(X^*)]^2, \end{aligned}$$

where  $(X^*, Y^*)$  is a copy of  $(X, Y)$  independent of the training data, and  $\mathbb{E}^*$  represents expectations taken over  $X^*$  and  $Y^*$  only. In particular, the rate of convergence of  $\mathcal{E}(\hat{\eta}_n)$  as the sample size  $n$  increases reflects the difficulty of the prediction problem. A closely related but different problem is that of estimating the intercept  $\alpha_0$  and the slope function  $\beta_0$ .

Many commonly used approaches to the prediction and estimation problems under the functional linear model (1) are based upon the functional principal component analysis (FPCA) (see, e.g., Ramsay and Silverman 2005; Yao, Müller, and Wang 2005; Cai and Hall 2006; Hall and Horowitz 2007). The functional principal components are determined solely by the observed functional predictors  $X_i$ . A crucial condition for the success of the FPCA-based methods is that the slope function  $\beta_0$  is efficiently represented in terms of the leading functional principal components. This condition, however, may not always be true. In many applications, it is not realistic to assume that the functional principal components form an efficient basis for the slope function  $\beta_0$  because the two are typically unrelated. When this condition fails to hold, the low-variance components of the predictor  $X$  have nonnegligible predictive power and the FPCA-based methods may not perform well. A similar phenomenon has also been observed in the performance of principal component regression (see, e.g., Jolliffe 1982) or singular value decomposition methods for linear inverse problems (see, e.g., Donoho 1995).

For illustration purposes, take the Canadian weather data as an example. It is one of the classic examples in functional linear regression and the goal is to predict the log annual precipitation at 35 different stations based on the average daily temperature. More detailed discussion of the data and analysis can be found in Ramsay and Silverman (2005) and Section 4 of the present article. The Fourier coefficients of the slope function with respect to the eigenfunctions of the sample covariance are estimated using the FPCA approach. The eigenvalues of the sample covariance and the estimated Fourier coefficients are shown in Figure 1. It is clear that although the eigenvalues decay nicely, the estimated Fourier coefficients do not decay at all. This is a typical example of the case when the slope function is not well represented by the leading principal components.

In this article, we study the minimax and adaptive prediction problems in the reproducing kernel Hilbert space (RKHS) framework under which the unknown slope function  $\beta_0$  is

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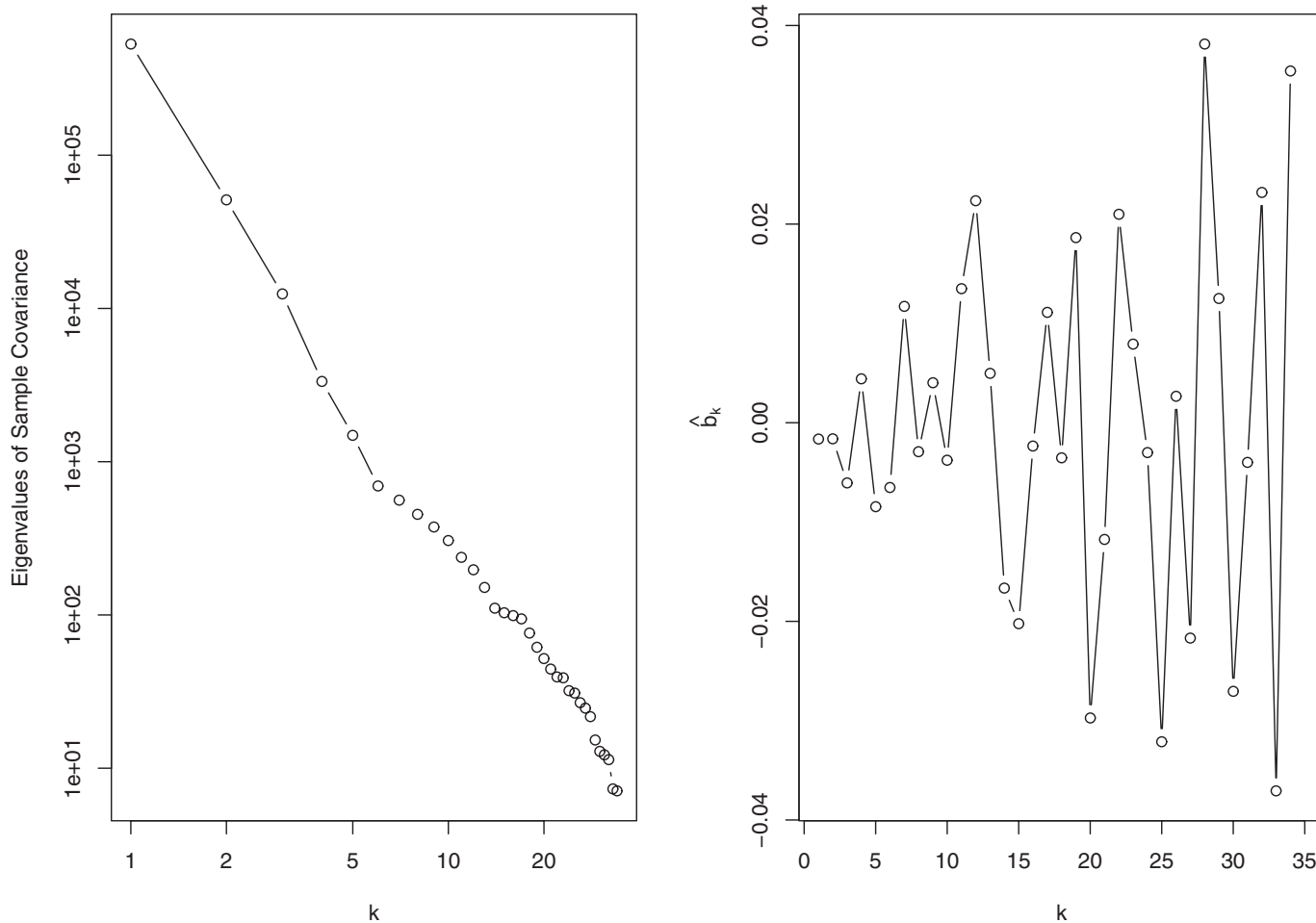


Figure 1. Canadian weather data: the eigenvalues of the sample covariance function are given in the left panel. Note that both axes in the left panel are in log scale. The right panel shows the estimated Fourier coefficients of the slope function with respect to the eigenfunctions of the sample covariance function.

assumed to reside in an RKHS  $\mathcal{H}(K)$  with a reproducing kernel  $K$ . The minimax rate of convergence of the excess prediction risk is established in a general setting with no constraint on the relationship between the reproducing kernel  $K$  and the covariance function  $C$  of the random predictor  $X$ . It is shown that under the functional linear model (1), the difficulty of the prediction problem as measured by the minimax rate of convergence depends on both kernels  $K$  and  $C$ . In particular, the alignment of  $K$  and  $C$  can significantly affect the optimal rate of convergence. The FPCA-based methods mentioned earlier correspond to the special setting where  $K$  and  $C$  are assumed to be perfectly aligned, that is,  $K$  and  $C$  share a common ordered set of eigenfunctions. The optimal rate of convergence of  $\mathcal{E}(\hat{\eta}_n)$  in this case is determined by the rate of decay of the product of the corresponding eigenvalues of  $K$  and  $C$  (Cai and Hall 2006). When  $K$  and  $C$  are not well aligned, as in the Canadian weather data example, the FPCA-based methods may not perform well.

The optimal rate of convergence for the prediction problem is established in two steps. First, a minimax lower bound is derived for the prediction problem. Then, a roughness regularization predictor is introduced and is shown to attain the rate of convergence given in the lower bound, when the tuning

parameter is appropriately chosen. This estimator is thus rate-optimal. The optimal choice of the tuning parameter depends, however, on the unknown covariance structure of the predictors  $X_i$ . A data-driven procedure for choosing the tuning parameter is then introduced. It is shown that the resulting procedure automatically achieves the optimal rate of convergence for a large collection of covariance functions. The adaptive procedure is easy to implement. Simulation studies are carried out to illustrate the merits of the adaptive predictor and to demonstrate the theoretical results.

The rest of the article is organized as follows. In Section 2, after basic notation and definitions are reviewed, we establish the optimal rate of convergence for the prediction problem by deriving both minimax lower and upper bounds. A roughness regularization predictor is shown to be rate-optimal when the tuning parameter is appropriately chosen. Section 3 proposes a data-driven method for choosing the tuning parameter and the resulting predictor is shown to adaptively achieve the optimal rate of convergence. Numerical experiments are reported in Section 4 to demonstrate the practical implications of the theoretical developments using both simulated and real datasets. The proofs are given in Section 6. We conclude with some discussions in Section 5.

## 2. OPTIMAL RATE OF CONVERGENCE

In this section, we establish the minimax rate of convergence of the excess prediction risk. The optimal rate is jointly determined by the reproducing kernel  $K$  and the covariance function  $C$ , and the alignment between  $K$  and  $C$  plays an important role. We begin by reviewing basic notation and properties regarding the reproducing kernel  $K$  and the covariance function  $C$ .

### 2.1 Notation and Definitions

Let  $\mathcal{T} \subset \mathbb{R}$  be a compact set. A reproducing kernel  $K : \mathcal{T} \times \mathcal{T} \rightarrow \mathbb{R}$  is a real, symmetric, square integrable, and nonnegative definite function. There is a one-to-one correspondence between a reproducing kernel  $K$  and an RKHS  $\mathcal{H}(K)$  which is a linear functional space endowed with an inner product  $\langle \cdot, \cdot \rangle_{\mathcal{H}(K)}$  such that for any  $t \in \mathcal{T}$ ,  $K(t, \cdot) \in \mathcal{H}(K)$ , and

$$f(t) = \langle K(t, \cdot), f \rangle_{\mathcal{H}(K)}, \quad \text{for any } f \in \mathcal{H}(K).$$

See, for example, Wahba (1990). Let  $X(\cdot)$  be a square integrable stochastic process defined over  $\mathcal{T}$ . The covariance function of  $X$  is also a real, symmetric, and nonnegative definite function defined as

$$C(s, t) = \mathbb{E}([X(s) - \mathbb{E}(X(s))][X(t) - \mathbb{E}(X(t))]), \quad \forall s, t \in \mathcal{T}.$$

The covariance function  $C$  is square integrable, if  $\mathbb{E}\|X\|_{\mathcal{L}_2}^2 < \infty$ , where

$$\|f\|_{\mathcal{L}_2}^2 = \langle f, f \rangle_{\mathcal{L}_2} \quad \text{and} \quad \langle f, g \rangle_{\mathcal{L}_2} = \int_{\mathcal{T}} f(t)g(t)dt.$$

For a real, symmetric, square integrable, and nonnegative definite function  $R : \mathcal{T} \times \mathcal{T} \rightarrow \mathbb{R}$ , let  $L_R : \mathcal{L}_2 \rightarrow \mathcal{L}_2$  denote an integral operator defined by

$$L_R(f)(\cdot) = \langle R(s, \cdot), f \rangle_{\mathcal{L}_2} = \int_{\mathcal{T}} R(s, \cdot)f(s)ds.$$

By the Riesz representation theorem,  $L_R$  can also be equivalently defined through

$$\langle f, L_R(g) \rangle_{\mathcal{H}(R)} = \langle f, g \rangle_{\mathcal{L}_2}.$$

The spectral theorem implies that there exists a set of orthonormalized eigenfunctions  $\{\psi_k^R : k \geq 1\}$  and a sequence of eigenvalues  $\theta_1^R \geq \theta_2^R \geq \dots > 0$  such that

$$R(s, t) = \sum_{k \geq 1} \theta_k^R \psi_k^R(s) \psi_k^R(t), \quad \forall s, t \in \mathcal{T},$$

and

$$L_R(\psi_k^R) = \theta_k^R \psi_k^R, \quad k = 1, 2, \dots$$

Throughout the article, we shall say that  $\{(\theta_k^R, \psi_k^R) : k \geq 1\}$  are the eigenvalue–eigenfunction pairs of the operator  $R$ , with the understanding that the pairs are ordered according to the eigenvalues in descending order,  $\theta_1^R \geq \theta_2^R \geq \dots$ . We say two linear operators (or their corresponding kernels) are perfectly aligned if they share the same ordered set of eigenfunctions (according to their eigenvalues in descending order).

Let  $L_{R^{1/2}}$  be a linear operator defined by

$$L_{R^{1/2}}(\psi_k^R) = \sqrt{\theta_k^R} \psi_k^R,$$

where

$$R^{1/2}(s, t) = \sum_{k \geq 1} \sqrt{\theta_k^R} \psi_k^R(s) \psi_k^R(t), \quad \forall s, t \in \mathcal{T}.$$

It is clear that  $L_{R^{1/2}} = L^{1/2}R$ . Moreover, define

$$(R_1 R_2)(s, t) = \int_{\mathcal{T}} R_1(s, u) R_2(u, t) dt.$$

Then,  $L_{R_1 R_2} = L_{R_1} \circ L_{R_2}$ .

For a given reproducing kernel  $K$  and a covariance function  $C$ , define the linear operator  $L_{K^{1/2}CK^{1/2}}$  by  $L_{K^{1/2}CK^{1/2}} = L_{K^{1/2}} \circ L_C \circ L_{K^{1/2}}$ , that is,

$$L_{K^{1/2}CK^{1/2}}(f) = L_{K^{1/2}}(L_C(L_{K^{1/2}}(f))).$$

If both  $L_{K^{1/2}}$  and  $L_C$  are bounded linear operators, so is  $L_{K^{1/2}CK^{1/2}}$ . By the spectral theorem, there exist a sequence of positive eigenvalues  $s_1 \geq s_2 \geq \dots > 0$  and a set of orthonormalized eigenfunctions  $\{\varphi_k : k \geq 1\}$  such that

$$K^{1/2}CK^{1/2}(s, t) = \sum_{k \geq 1} s_k \varphi_k(s) \varphi_k(t), \quad \forall s, t \in \mathcal{T},$$

and

$$L_{K^{1/2}CK^{1/2}}(\varphi_k) = s_k \varphi_k, \quad k = 1, 2, \dots$$

It is easy to see that the eigenvalues  $\{s_k : k \geq 1\}$  of the linear operator  $L_{K^{1/2}CK^{1/2}}$  depend on the eigenvalues of both the reproducing kernel  $K$  and the covariance function  $C$  as well as the alignment between  $K$  and  $C$ . We shall show in Sections 2.2 and 2.3 that the difficulty of the prediction problem as measured by the minimax rate of convergence of the excess prediction risk is determined by the decay rate of the eigenvalues  $\{s_k : k \geq 1\}$ .

For two sequences  $\{a_k : k \geq 1\}$  and  $\{b_k : k \geq 1\}$  of positive real numbers, we write  $a_k \asymp b_k$ , if there are positive constants  $c$  and  $C$  independent of  $k$  such that  $c \leq a_k/b_k \leq C$  for all  $k \geq 1$ .

### 2.2 Minimax Lower Bound

The optimal rate of convergence of the excess prediction risk will be established in two steps. We first derive a minimax lower bound in this section and then show in Section 2.3 the convergence rate of the lower bound is in fact optimal by constructing a predictor that attains this rate of convergence. The minimax lower bound is given in the following theorem.

*Theorem 1.* Suppose the eigenvalues  $\{s_k : k \geq 1\}$  of the linear operator  $L_{K^{1/2}CK^{1/2}}$  satisfy  $s_k \asymp k^{-2r}$  for some constant  $0 < r < \infty$ , then the excess prediction risk satisfies

$$\lim_{a \rightarrow 0} \liminf_{n \rightarrow \infty} \sup_{\tilde{\eta}} \mathbb{P} \left\{ \mathcal{E}(\tilde{\eta}) \geq an^{-\frac{2r}{2r+1}} \right\} = 1, \quad (2)$$

where the infimum is taken over all possible predictors  $\tilde{\eta}$  based on the training data  $\{(X_i, Y_i) : i = 1, \dots, n\}$ .

It is interesting to compare Theorem 1 with some of the known results in the literature in which the reproducing kernel  $K$  and the covariance function  $C$  are assumed to be perfectly aligned, that is, they share the same ordered set of eigenfunctions. The minimax lower bound obtained here generalizes the

earlier results for this special case. Let

$$X(t) = \sum_{k=1}^{\infty} Z_k \psi_k(t)$$

be the Karhunen–Loève decomposition of  $X$ , where  $\{Z_k : k \geq 1\}$  are uncorrelated random variables and  $\{\psi_k : k \geq 1\}$  is an orthonormal basis of  $\mathcal{L}_2(\mathcal{T})$ . Set  $\theta_k^C = \text{var}(Z_k)$ . We shall assume that  $\theta_k^C$  are indexed in descending order,  $\theta_1^C \geq \theta_2^C \geq \dots$ . Then, it is clear that  $\{(\theta_k^C, \psi_k) : k \geq 1\}$  also constitutes the eigenvalue–eigenfunction pairs of the covariance function  $C$ , that is,

$$C(s, t) = \sum_{k=1}^{\infty} \theta_k^C \psi_k(s) \psi_k(t), \quad \text{for } s, t \in \mathcal{T}.$$

Consider the case where the reproducing kernel  $K$  is perfectly aligned with  $C$ , that is,

$$K(s, t) = \sum_{k=1}^{\infty} \theta_k^K \psi_k(s) \psi_k(t), \quad \text{for } s, t \in \mathcal{T},$$

with  $\theta_1^K \geq \theta_2^K \geq \dots \geq 0$  being the eigenvalues of  $K$ . In this case, it is easy to see that

$$L_{K^{1/2}CK^{1/2}}(\psi_k) = \theta_k^K \theta_k^C \psi_k, \quad k = 1, 2, \dots,$$

indicating that the eigenvalues  $s_k = \theta_k^K \theta_k^C$ ,  $k = 1, 2, \dots$ , and  $s_1 \geq s_2 \geq \dots$ . If  $\theta_k^K \asymp k^{-2r_1}$  and  $\theta_k^C \asymp k^{-2r_2}$ , then  $s_k$  decays at the rate  $k^{-2(r_1+r_2)}$  and by Theorem 1,

$$\lim_{a \rightarrow 0} \liminf_{n \rightarrow \infty} \sup_{\tilde{\eta}} \mathbb{P} \left\{ \mathcal{E}(\tilde{\eta}) \geq an^{-\frac{2(r_1+r_2)}{2(r_1+r_2)+1}} \right\} = 1.$$

This special setting coincides with those considered in Cai and Hall (2006) and Yuan and Cai (2010). Similar results have been established earlier in these articles for this special setting.

In general, however, the eigenvalues of  $K$  and  $C$  alone cannot determine the decay rate of the eigenvalues of  $L_{K^{1/2}CK^{1/2}}$ . For example, when  $\theta_k^K \asymp k^{-2r_1}$ ,  $\theta_k^C \asymp k^{-2r_2}$ , and  $\psi_k^C = \psi_{k^2}^K$  for  $k = 1, 2, \dots$ , then  $s_k \asymp k^{-(4r_1+2r_2)}$ .

### 2.3 Minimax Upper Bound

Section 2.2 developed a minimax lower bound for the excess prediction risk. We shall now consider the minimax upper bound and show that the lower bound established in Theorem 1 can in fact be achieved. We shall construct a predictor using a roughness regularization method and show that the predictor is asymptotically rate-optimal.

One of the most commonly used methods in nonparametric function estimation is the roughness regularization method (see, e.g., Ramsay and Silverman 2005) where the intercept  $\alpha_0$  and the slope function  $\beta_0$  are estimated by

$$(\hat{\alpha}_{n\lambda}, \hat{\beta}_{n\lambda}) = \underset{a \in \mathbb{R}, b \in \mathcal{H}(K)}{\operatorname{argmin}} \left\{ \sum_{i=1}^n \left( Y_i - a - \int_{\mathcal{T}} X_i(t) b(t) dt \right)^2 + \lambda \|b\|_{\mathcal{H}(K)}^2 \right\}. \quad (3)$$

Here,  $\lambda > 0$  is a tuning parameter that balances the tradeoff between the fidelity to the data measured by the sum of squares and the smoothness of the estimate measured by the squared RKHS norm. The estimate  $\hat{\beta}_\lambda$  is readily computable even though

the minimization in Equation (3) is taken over an infinitely dimensional function space  $\mathcal{H}(K)$ . More specifically,  $\hat{\beta}_{n\lambda}$  can be expressed as

$$\hat{\beta}_{n\lambda}(\cdot) = \sum_{i=1}^n c_i \int_{\mathcal{T}} K(t, \cdot) X_i(t) dt \quad (4)$$

for some  $c_1, \dots, c_n \in \mathbb{R}$ , and they can be computed together with  $\hat{\alpha}_{n\lambda}$  by plugging (4) back to (3). The readers are referred to Yuan and Cai (2010) for more details on the implementation.

Given the estimates  $\hat{\alpha}_{n\lambda}$  and  $\hat{\beta}_{n\lambda}$ , the predictor  $\hat{\eta}_{n\lambda}$  is obtained by

$$\hat{\eta}_{n\lambda}(X) = \hat{\alpha}_{n\lambda} + \int_{\mathcal{T}} X(t) \hat{\beta}_{n\lambda}(t) dt.$$

The following theorem states that with an appropriately chosen  $\lambda$ , the predictor  $\hat{\eta}_{n\lambda}$  attains the convergence rate of the lower bound given by Theorem 1 and is therefore rate-optimal.

*Theorem 2.* Assume that there exists a constant  $c > 0$  such that for any square integrable function  $f$  defined over the domain  $\mathcal{T}$ ,

$$\mathbb{E} \left( \int_{\mathcal{T}} X(t) f(t) dt \right)^4 \leq c \left( \mathbb{E} \left( \int_{\mathcal{T}} X(t) f(t) dt \right)^2 \right)^2. \quad (5)$$

Suppose the eigenvalues  $\{s_k : k \geq 1\}$  satisfy  $s_k \asymp k^{-2r}$  for some constant  $0 < r < \infty$ , then

$$\lim_{A \rightarrow \infty} \liminf_{n \rightarrow \infty} \sup_{\beta_0 \in \mathcal{H}(K)} \mathbb{P} \left\{ \mathcal{E}(\hat{\eta}_{n\lambda}) \geq An^{-\frac{2r}{2r+1}} \right\} = 0, \quad (6)$$

provided that  $\lambda \asymp n^{-2r/(2r+1)}$ .

Condition (5) states that linear functionals of  $X$  have bounded kurtosis that is satisfied in particular with  $c = 3$  when  $X$  follows a Gaussian process.

Theorems 1 and 2 together show that the minimax rate of convergence for the excess prediction risk is

$$n^{-2r/(2r+1)},$$

which is determined by the rate of decay of the eigenvalues of the operator  $L_{K^{1/2}CK^{1/2}}$ . The optimal rate of convergence depends not only on the eigenvalues of  $K$  and  $C$  but also on how their eigenfunctions align with each other.

### 3. ADAPTIVE PREDICTION

Section 2 established the minimax rate of convergence for the excess prediction risk. As shown in Theorem 2, the optimal rate can be attained by the roughness regularization predictor  $\hat{\eta}_{n\lambda}$  when the tuning parameter  $\lambda$  is chosen appropriately. However, the proper choice of  $\lambda$  depends on  $r$ , which is unknown since it is determined by the linear operator  $L_{K^{1/2}CK^{1/2}}$  and the covariance function  $C$  is not known a priori. It is important to develop a data-driven choice of  $\lambda$  that does not require the knowledge of  $C$ . In this section, we introduce such an adaptive method for choosing  $\lambda$ .

To motivate our procedure, recall that  $r$  represents the decay rate of the eigenvalues of  $L_{K^{1/2}CK^{1/2}}$ , which can be naturally

approximated by  $L_{K^{1/2}C_n K^{1/2}}$ , where

$$C_n(s, t) = \frac{1}{n} \sum_{i=1}^n (X_i(s) - \bar{X}(s))(X_i(t) - \bar{X}(t)).$$

Observe that

$$\begin{aligned} \langle f, L_{K^{1/2}C_n K^{1/2}} g \rangle_{\mathcal{L}_2} &= \langle L_{K^{1/2}} f, L_{C_n} L_{K^{1/2}} g \rangle_{\mathcal{L}_2} \\ &= \frac{1}{n} \sum_{i=1}^n \langle L_{K^{1/2}} f, X_i - \bar{X} \rangle_{\mathcal{L}_2} \\ &\quad \times \langle L_{K^{1/2}} g, X_i - \bar{X} \rangle_{\mathcal{L}_2} \\ &= \frac{1}{n} \sum_{i=1}^n \langle f, L_{K^{1/2}} X_i - \bar{X} \rangle_{\mathcal{L}_2} \\ &\quad \times \langle g, L_{K^{1/2}} X_i - \bar{X} \rangle_{\mathcal{L}_2} \\ &= \langle f, L_{H_n} g \rangle_{\mathcal{L}_2}, \end{aligned}$$

where

$$H_n(s, t) = \frac{1}{n} \sum_{i=1}^n (L_{K^{1/2}}(X_i - \bar{X}))(s)(L_{K^{1/2}}(X_i - \bar{X}))(t).$$

By duality, the eigenvalues of  $H_n$  are also the eigenvalues of the Gram matrix  $G = (G_{ij})_{1 \leq i, j \leq n}$ , where

$$\begin{aligned} G_{ij} &= \frac{1}{n} \langle L_{K^{1/2}}(X_i - \bar{X}), L_{K^{1/2}}(X_j - \bar{X}) \rangle_{\mathcal{L}_2} \\ &= \frac{1}{n} \int_{\mathcal{T}^2} (X_i - \bar{X})(s)K(s, t)(X_j - \bar{X})(t)dsdt. \end{aligned}$$

Information on  $r$  can thus be recovered from the eigenvalues of  $G$ .

Write

$$\gamma_n(\delta) = \left( \frac{1}{n} \sum_{k \geq 1} \min\{s_k, \delta^2\} \right)^{1/2}$$

and define

$$\rho(C, K) := \inf\{\rho \geq n^{-1} \log n : \gamma_n(\delta) \leq \rho^{1/2} \delta + \rho, \forall \delta \in [0, 1]\}.$$

It is not hard to see that  $\rho \asymp n^{-2r/(2r+1)}$ , if  $s_k \asymp k^{-2r}$  (see, e.g., Mendelson 2002). Therefore, we can use an estimate of  $\rho$  as our choice of the tuning parameter in defining  $\hat{\eta}_\lambda$ . To this end, we consider the sample analog of  $\rho$ . Denote by  $\hat{s}_1 \geq \hat{s}_2 \geq \dots \geq \hat{s}_n$  the eigenvalues of  $G$ . Write

$$\hat{\rho}(G) := \inf\{\rho \geq n^{-1} \log n : \hat{\gamma}_n(\delta) \leq \rho^{1/2} \delta + \rho, \forall \delta \in [0, 1]\},$$

where

$$\hat{\gamma}_n(\delta) = \left( \frac{1}{n} \sum_{k=1}^n \min\{\hat{s}_k, \delta^2\} \right)^{1/2}.$$

Theorem 3 shows that with high probability  $\hat{\rho}$  is of the same order as  $\rho$  whenever  $\|L_{K^{1/2}} X\|_{\mathcal{L}_2}$  has exponential tails.

*Theorem 3.* Assume that there exist some constants  $c_1, c_2 > 0$  such that

$$\mathbb{P}\{\|L_{K^{1/2}} X\|_{\mathcal{L}_2} \geq x\} \leq c_1 \exp(-c_2 x^2), \quad \text{for } x \geq 0. \quad (7)$$

Then, there are constants  $0 < c_3 < c_4 < \infty$  such that

$$\mathbb{P}\left\{c_3 \leq \frac{\hat{\rho}}{\rho} \leq c_4\right\} \rightarrow 1$$

as  $n \rightarrow \infty$ .

We note that the tail condition (7) for  $\|L_{K^{1/2}} X\|_{\mathcal{L}_2}$  can also be replaced by

$$\mathbb{P}\{\|X\|_{\mathcal{L}_2} \geq x\} \leq c_1 \exp(-c_2 x^2), \quad \text{for } x \geq 0,$$

because  $\|L_{K^{1/2}} X\|_{\mathcal{L}_2} \leq \theta_1^K \|X\|_{\mathcal{L}_2}$ . It holds true, in particular, if  $X$  is a Gaussian process.

The following result shows that the predictor  $\hat{\eta}_{n\lambda}$  with the data-driven choice of the tuning parameter  $\lambda$  indeed adaptively achieves the optimal rate of convergence.

*Theorem 4.* Assume that there exist constants  $c_1, c_2 > 0$  such that

$$\mathbb{P}\{\|L_{K^{1/2}} X\|_{\mathcal{L}_2} \geq x\} \leq c_1 \exp(-c_2 x^2), \quad \text{for } x \geq 0,$$

and for any  $f \in \mathcal{L}_2$

$$\mathbb{E}\left(\int_{\mathcal{T}} X(t)f(t)dt\right)^4 \leq c \left(\mathbb{E}\left(\int_{\mathcal{T}} X(t)f(t)dt\right)^2\right)^2. \quad (8)$$

Suppose the eigenvalues  $\{s_k : k \geq 1\}$  satisfy  $s_k \asymp k^{-2r}$  for some constant  $0 < r < \infty$ , then

$$\lim_{A \rightarrow \infty} \lim_{n \rightarrow \infty} \sup_{\beta_0 \in \mathcal{H}(K)} \mathbb{P}\left\{\mathcal{E}(\hat{\eta}_{n\lambda}) \geq A n^{-\frac{2r}{2r+1}}\right\} = 0 \quad (9)$$

for any  $\lambda \asymp \hat{\rho}$ .

In light of Theorem 4, we can choose  $\lambda = \hat{\lambda} := c\hat{\rho}$  for some constant  $c > 0$  to ensure that the tuning parameter is of the appropriate order. In practice, the constant  $c$  can be further fine-tuned to ensure enhanced performance. For example, our experience suggests that  $\hat{\lambda} = \hat{\rho}/\hat{s}_1^2$  usually leads to satisfactory performance.

## 4. NUMERICAL EXPERIMENTS

We turn in this section to the numerical performance of the proposed adaptive predictor and demonstrate the practical implications of the theoretical results developed in the last two sections. We shall begin with simulation results, and then return to the analysis of the Canadian weather data mentioned in the introduction.

### 4.1 Simulation Study

To fix our ideas, we shall focus on the case where  $\mathcal{T} = [0, 1]$  and  $\mathcal{H}(K)$  consists of functions in the linear span of the cosine basis,

$$f(t) = \sqrt{2} \sum_{k \geq 1} f_k \cos(k\pi t), \quad t \in [0, 1],$$

such that

$$\sum_{k \geq 1} k^4 f_k^2 < \infty.$$

When endowed with norm

$$\begin{aligned} \|f\|_{\mathcal{H}(K)}^2 &= \int (f'')^2 = \int_0^1 \left( \sqrt{2} \sum_{k \geq 1} (k\pi)^2 f_k \cos(k\pi t) \right)^2 \\ &= \sum_{k \geq 1} (k\pi)^4 f_k^2, \end{aligned}$$

$\mathcal{H}(K)$  forms an RKHS with the reproducing kernel

$$\begin{aligned} K(s, t) &= \sum_{k \geq 1} \frac{2}{(k\pi)^4} \cos(k\pi s) \cos(k\pi t) \\ &= \sum_{k \geq 1} \frac{1}{(k\pi)^4} \cos(k\pi(s-t)) + \sum_{k \geq 1} \frac{1}{(k\pi)^4} \cos(k\pi(s+t)) \\ &= -\frac{1}{3} (B_4(|s-t|/2) + B_4((s+t)/2)), \end{aligned}$$

where  $B_k$  is the  $k$ th Bernoulli polynomial. Here, the following fact (Abramowitz and Stegun 1965) is used:

$$B_{2m}(x) = (-1)^{m-1} 2(2m)! \sum_{k \geq 1} \frac{\cos(2\pi kx)}{(2\pi k)^{2m}}, \quad \forall x \in [0, 1].$$

In this setting, the roughness regularization estimate defined by (3) is given by

$$\begin{aligned} (\hat{\alpha}_{n\lambda}, \hat{\beta}_{n\lambda}) &= \operatorname{argmin}_{a \in \mathbb{R}, b \in \mathcal{H}(K)} \left\{ \sum_{i=1}^n \left( Y_i - a - \int_{\mathcal{T}} X_i(t) b(t) dt \right)^2 \right. \\ &\quad \left. + \lambda \int (b'')^2 \right\}. \end{aligned}$$

For brevity, we shall also assume that there is no intercept in the functional linear model (1). The roughness regularization estimate then becomes

$$\hat{\beta}_{n\lambda} = \operatorname{argmin}_{b \in \mathcal{H}(K)} \left\{ \sum_{i=1}^n \left( Y_i - \int_{\mathcal{T}} X_i(t) b(t) dt \right)^2 + \lambda \int (b'')^2 \right\}.$$

To investigate the effect of varying covariance function, the true slope function is taken to be

$$\beta_0(t) = \sum_{k \geq 1} 4\sqrt{2}(-1)^{k-1} k^{-2} \cos(k\pi t).$$

We first consider the effect of varying smoothness of the covariance function. More specifically, set

$$C(s, t) = \sum_{k \geq 1} 2k^{-2r_2} \cos(k\pi s) \cos(k\pi t).$$

The parameter  $r_2$  controls how fast the eigenvalues of the covariance function  $C$  decay and therefore by Theorem 1 determines the optimal rates of convergence. We let the value of  $r_2$  vary between 1 and 3. The functional predictors were simulated from a centered Gaussian process with covariance function  $C$ . To comprehend the trend as the sample size increases, the sample size  $n$  is chosen to be  $n = 32, 64, 128, 256, 512, \text{ and } 1024$ . For each sample size and each value of  $r_2$ , we simulate data from the functional linear model with  $\sigma = 0.5$ . Both the roughness regularization and functional principal component estimates were computed and the tuning parameters,  $\lambda$  for the former and the number of principal components for the latter, were chosen to yield the smallest excess risk so that it reflects the optimal rate achieved by both methods. The experiment was repeated for 200 times and the results are summarized in Figure 2.

In this example, both the reproducing kernel and the covariance function share a common ordered set of eigenfunctions. In this case, as shown previously by Cai and Hall (2006) and Yuan and Cai (2010), both methods can attain the optimal rate of convergence. The similarity between the two methods as observed in Figure 2 simply confirms these earlier findings.

One of the key messages from Theorem 1 is that in addition to the decay of the eigenvalues of the reproducing kernel and covariance function, the alignment of the eigenfunctions of the two kernels also plays a crucial role in determining the optimal rate of convergence. To demonstrate the effect of such an alignment, we now consider a slightly different covariance function:

$$C(s, t) = \sum_{k \geq 1} 2\theta_k \cos(k\pi s) \cos(k\pi t),$$

where

$$\theta_k = (|k - k_0| + 1)^{-2}.$$

In this setting, the leading eigenfunctions of  $C$  are located around the  $k_0$ th eigenfunction of the reproducing kernel and in a certain sense controls the misalignment between the covariance function and the reproducing kernel. We consider  $k_0 = 5, 10, 15, \text{ and } 20$  to appreciate the effect of the alignment. The simulation was carried out in a similar fashion as before and the results are reported in the top panels of Figure 3.

As expected, for larger values of  $k_0$ , the functional principal component-based approach performs rather poorly when the sample size is small. However, as one can observe from Figure 3, both the methods appear to converge at similar rates as the sample size increases although the roughness regularization method performs better. The top right panel displays the median relative efficiency of the roughness regularization over the functional principal component-based method defined as  $\mathcal{E}(\hat{\eta}^{\text{FPCA}})/\mathcal{E}(\hat{\eta}^{\text{REG}})$ , where  $\hat{\eta}^{\text{REG}}$  and  $\hat{\eta}^{\text{FPCA}}$  represent the two estimates, respectively. It is evident that the efficiency of the roughness regularization increases with  $k_0$ , which reflects how poorly the eigenfunctions of  $K$  and  $C$  align. In most cases, the roughness regularization estimate outperforms the functional principal component-based method by an order of magnitude. It is noteworthy that the performance of FPCA-based approach quickly deteriorates as  $k_0$  increases. In particular, when  $k_0$  is greater than the sample size, the FPCA approach will fail since the true slope function is orthogonal to the leading functional principal components. In this example, we have chosen the largest value of  $k_0$  to be 20 to ensure that  $k_0$  is always smaller than the sample size  $n$ , and to give a better contrast of the two methods.

To further demonstrate the effect of alignment between the eigenfunctions of  $K$  and  $C$ , we also considered a different set of covariance functions. In particular,

$$C(s, t) = \sum_{k \geq 1} 2k^{-2r_2} h_k(s) h_k(t),$$

where the  $h_k$ 's are the Haar functions, that is,

$$h_{2^k+l-1}(t) = \begin{cases} 2^{k/2}, & t \in \left[ \frac{l-1}{2^k}, \frac{l-1/2}{2^k} \right), \\ -2^{k/2}, & t \in \left[ \frac{l-1/2}{2^k}, \frac{l}{2^k} \right), \\ 0, & \text{otherwise,} \end{cases}$$

for  $k = 0, 1, 2, \dots$  and  $l = 1, \dots$ . Different from the previous two simulation settings, the eigenfunctions of  $C$  are the Haar basis of  $\mathcal{L}_2$  that are different from those of  $K$ . We again apply both the roughness regularization and functional principal component-based methods to each simulated dataset and

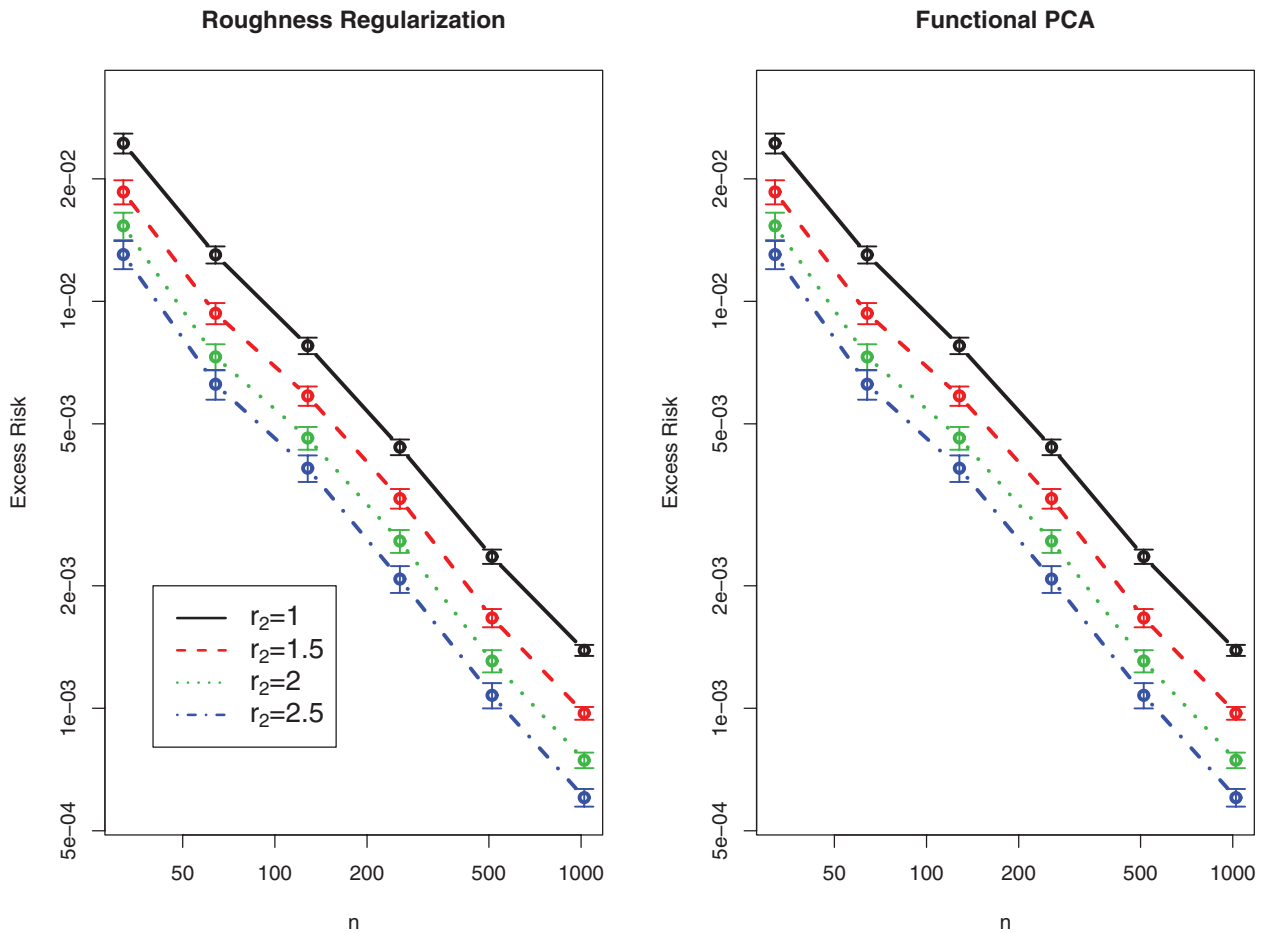


Figure 2. Effect of smoothness of the covariance function: two hundred datasets were simulated for each combination of sample size  $n = 32, 64, 128, 256, 512,$  or  $1024$  and  $r_2 = 1, 1.5, 2, 2.5,$  or  $3$ . Both the roughness regularization estimate and the functional principal component estimate were computed with tuning parameter chosen to minimize the integrated squared error. The circle corresponds to the excess risk averaged over 200 datasets and error bars correspond to the mean  $\pm 1$  SE. Both axes are in log scale. The online version of this figure is in color.

summarize the findings in the lower panels of Figure 3. Similar to before, the results are averaged over 200 simulated datasets for each value of  $r_2$ . We observe similar comparison as in the last example with the roughness regularization significantly outperforming the functional principal component-based method.

We now turn to the choice of the tuning parameter  $\lambda_n$ . The adaptive prediction procedure is applied to the simulated data. For brevity, we focus on the first two simulation settings. For each simulated dataset, the tuning parameter  $\lambda$  is chosen as  $\lambda = \hat{\rho}/\hat{s}_1^2$ . As mentioned earlier, this particular scaling usually yields fairly reasonable results, although it is plausible that it can be further improved by using other data-driven choice of scaling factors. Figure 4 summarizes the excess risk for all settings. The behavior of the resulting estimate closely resembles those reported earlier, indicating that the adaptive procedure proposed earlier is indeed capable of achieving the optimal rate of convergence.

#### 4.2 Canadian Weather Data

Finally, we revisit the Canadian weather data example. The dataset, one of the most popular examples in functional linear regression, contains daily temperature and precipitation at 35 different locations in Canada averaged over 1960–1994. The goal is to predict the log annual precipitation based on the aver-

age daily temperature. As shown in Figure 1, the application of functional PCA-based methods could be problematic since the estimated slope function does not seem to allow for a compact representation with respect to the eigenfunctions of the covariance function. Alternatively, we applied the regularization approach discussed earlier to the data. The data were collected on the basis of calendar year. The nature of the data suggests that the predictor function  $X$  be periodic on  $[0, 1]$  (year), and so is the slope function  $\beta_0$ . Let  $\mathcal{W}_2^{\text{per}}$  be the second-order Sobolev space of periodic functions on  $[0, 1]$ , we estimate  $\alpha_0$  and  $\beta_0$  by

$$(\hat{\alpha}, \hat{\beta}) = \underset{a \in \mathbb{R}, b \in \mathcal{W}_2^{\text{per}}}{\text{argmin}} \left\{ \left[ Y_i - a - \int_0^1 X_i(t)b(t)dt \right]^2 + \lambda \int_0^1 [b''(t)]^2 dt \right\}. \tag{10}$$

The space  $\mathcal{W}_2^{\text{per}}$ , endowed with the norm

$$\|b\|_{\mathcal{W}_2^{\text{per}}}^2 = \left[ \int_0^1 b(t)dt \right]^2 + \int_0^1 [b''(t)]^2 dt,$$

has a reproducing kernel:

$$K(s, t) = 1 - \frac{1}{24} B_4(|s - t|), \quad \forall s, t \in [0, 1],$$

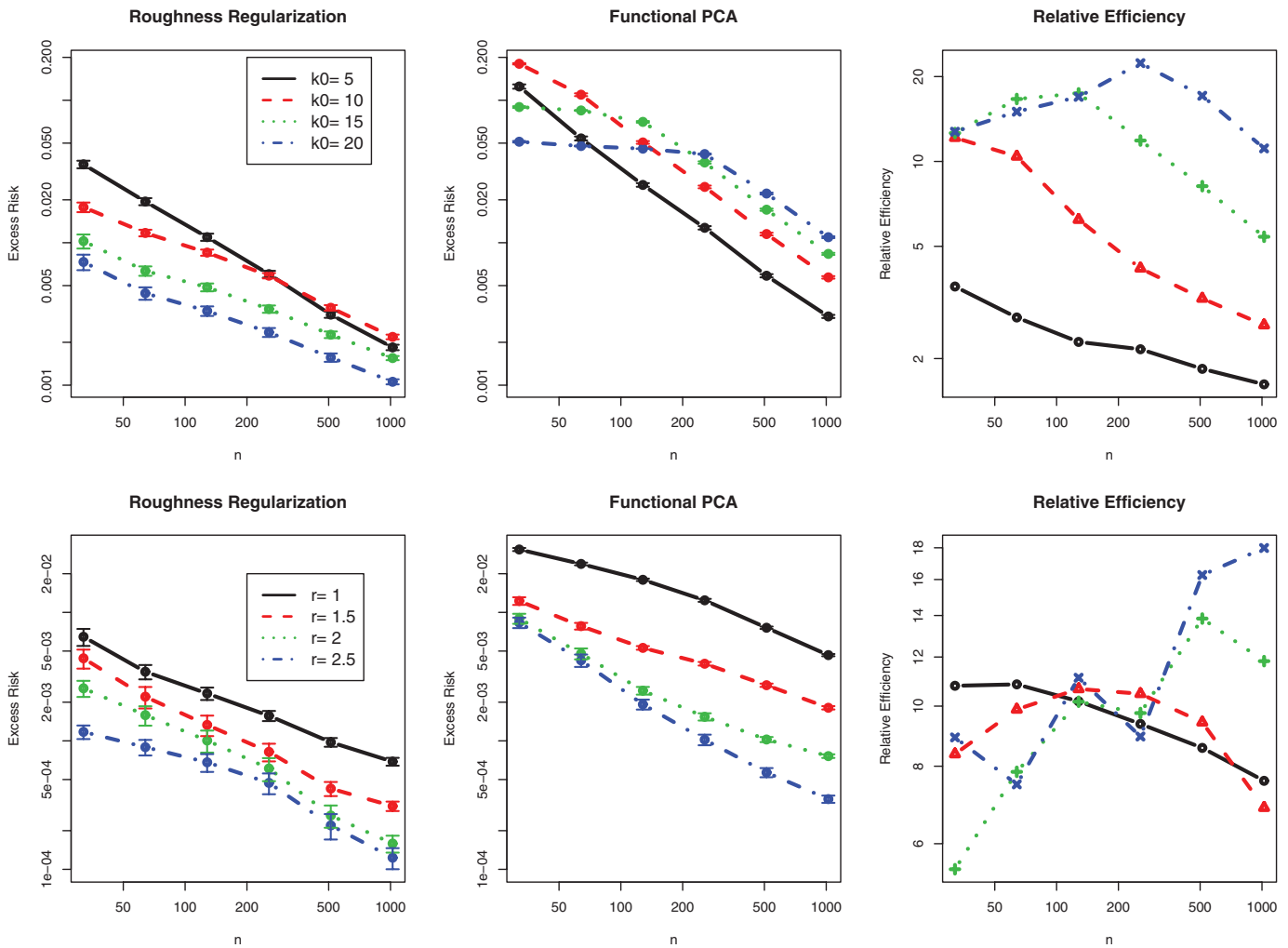


Figure 3. Effect of alignment of eigenfunctions between the reproducing kernel and the covariance function: for the top panels, 200 datasets were simulated for each combination of sample size  $n = 32, 64, 128, 256, 512,$  or  $1024$  and the location of the first principal component  $k_0 = 5, 10, 15,$  or  $20$ . For the bottom panels, 200 datasets were simulated for each combination of sample size  $n = 32, 64, 128, 256, 512,$  or  $1024$  and the decay rate of the eigenvalues of  $C, r_2 = 1, 1.5, 2, 2.5$ . In the left and middle panels, the circle corresponds to the excess risk averaged over 200 datasets and error bars correspond to the mean  $\pm 1$  SE. In the rightmost panels, median relative efficiency of the roughness regularization method over functional principal component-based method is given for the two simulation settings. Both axes are in log scale. The online version of this figure is in color.

where  $B_4$  is the fourth Bernoulli polynomial (see, e.g., Wahba 1990). As shown in Yuan and Cai (2010), the solution of Equation (10) can be expressed as

$$\hat{\alpha} = \bar{Y} - \int_0^1 \bar{X}(t)\hat{\beta}(t)dt$$

and

$$\hat{\beta}(\cdot) = c_0 + \sum_{i=1}^n c_i \int_0^1 (X_i - \bar{X})(s)K_1(s, \cdot)ds,$$

where

$$K_1(s, t) = -\frac{1}{24}B_4(|s - t|)$$

is the reproducing kernel of the orthogonal complement of constant functions in  $\mathcal{W}_2^{\text{per}}$ . With this representation, the objective function of (10) becomes quadratic in terms of  $c_0, c_1, \dots, c_n$  and they can be solved through

$$(c_0, (c_1, \dots, c_n)^T) = \underset{a_0 \in \mathbb{R}, \mathbf{a} \in \mathbb{R}^n}{\text{argmin}} \{ \|\mathbf{Y} - \Sigma \mathbf{a} - a_0 \mathbf{u}\|^2 + \lambda \mathbf{a}^T \Sigma \mathbf{a} \},$$

where  $\mathbf{Y} = (Y_1, \dots, Y_n)^T, \Sigma$  is an  $n \times n$  matrix whose  $(i, j)$  entry is

$$\Sigma_{ij} = \int_0^1 (X_i - \bar{X})(s)K_1(s, t)(X_j - \bar{X})(t)dsdt,$$

and  $\mathbf{u}$  is an  $n$ -dimensional vector with the  $i$ th entry given by

$$u_i = \int_0^1 (X_i - \bar{X})(s)ds.$$

The integrals in defining  $\Sigma_{ij}$  and  $u_i$  can be approximated by summations for practical purposes. We also used the proposed adaptive procedure to choose the tuning parameter. The estimated slope function is given in the left panel of Figure 5. The right panel provides the normal Q-Q plot of the residuals, which suggests a fairly good fit of the data. The plot suggests that the precipitation at Kamloops station, corresponding to the rightmost point in the Q-Q plot, may be overestimated using the functional linear model. This can be attributed to the fact that Kamloops lies deep in the Thompson River Valley. These



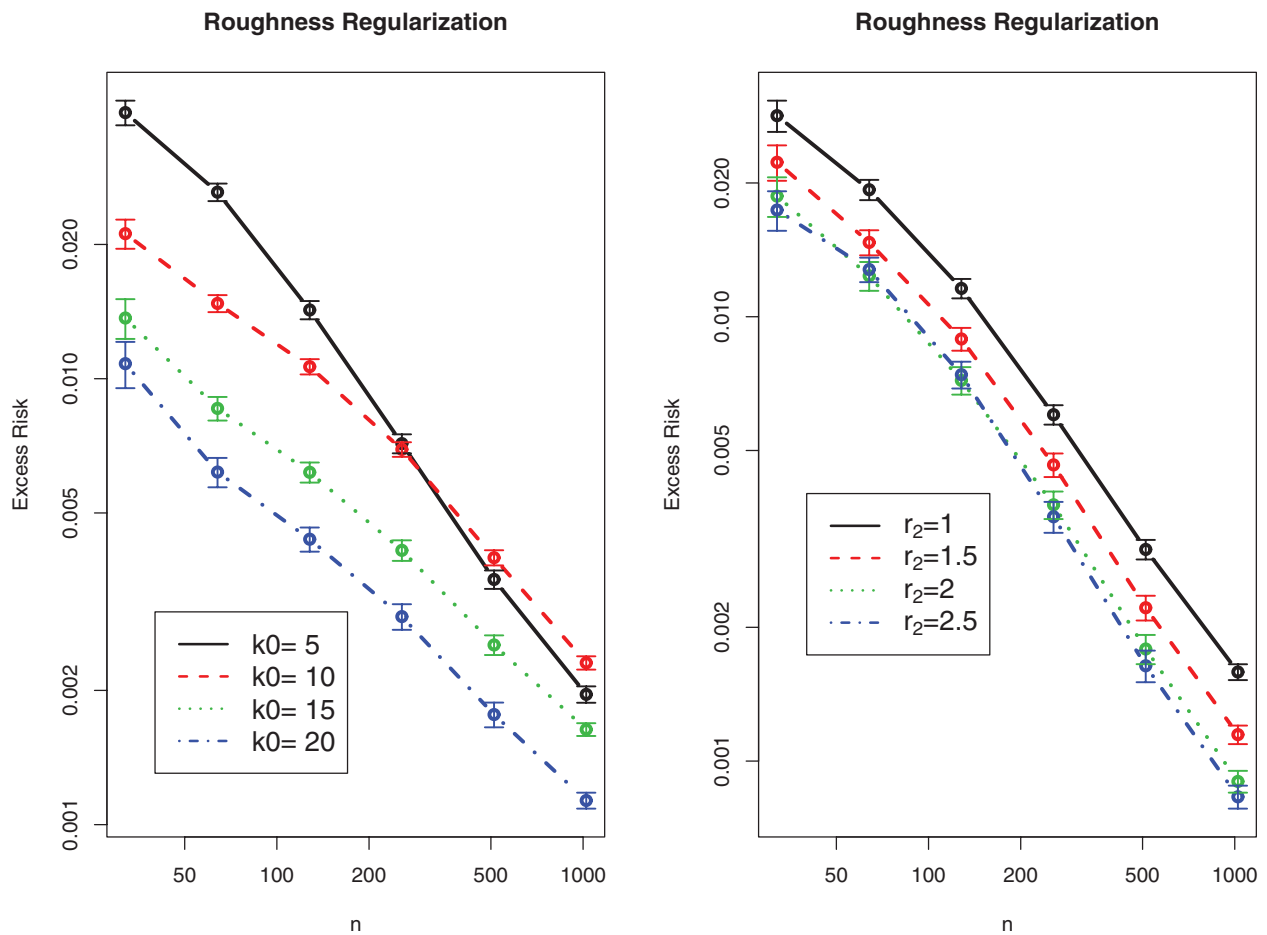


Figure 4. Excess risk of the adaptive prediction: each model used before was considered and the tuning parameter is chosen to be  $\hat{\rho}/\hat{s}_1^2$ . The left panel shows the result for different combinations of  $n$  and  $k_0$  whereas the right panel corresponds to the result for different combinations of  $n$  and  $r_0$ . The errors bars correspond to the mean  $\pm 1$  SE. Both axes are in log scale. The online version of this figure is in color.

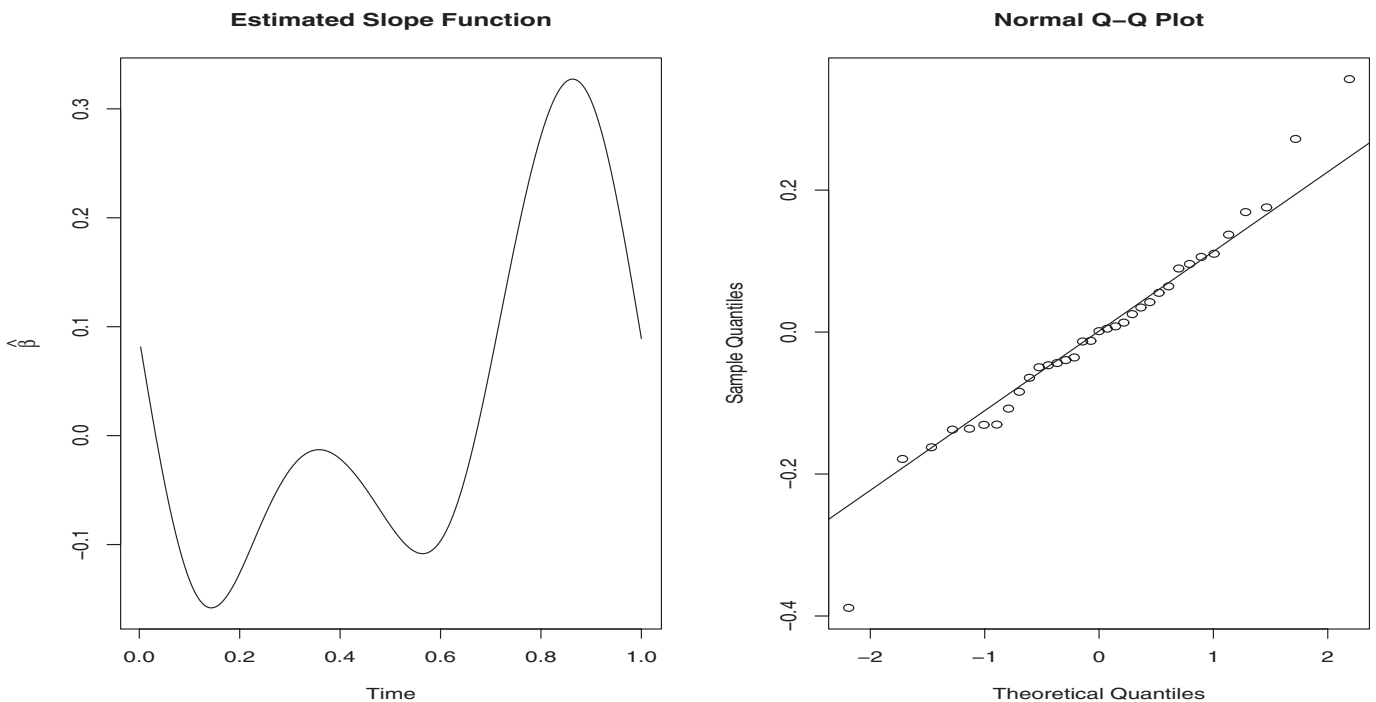


Figure 5. Canadian weather data: smoothness regularization estimate of the slope function with adaptive choice of the tuning parameter. The left panel gives the estimated slope function whereas the right panel shows the Q-Q plot of the residuals.

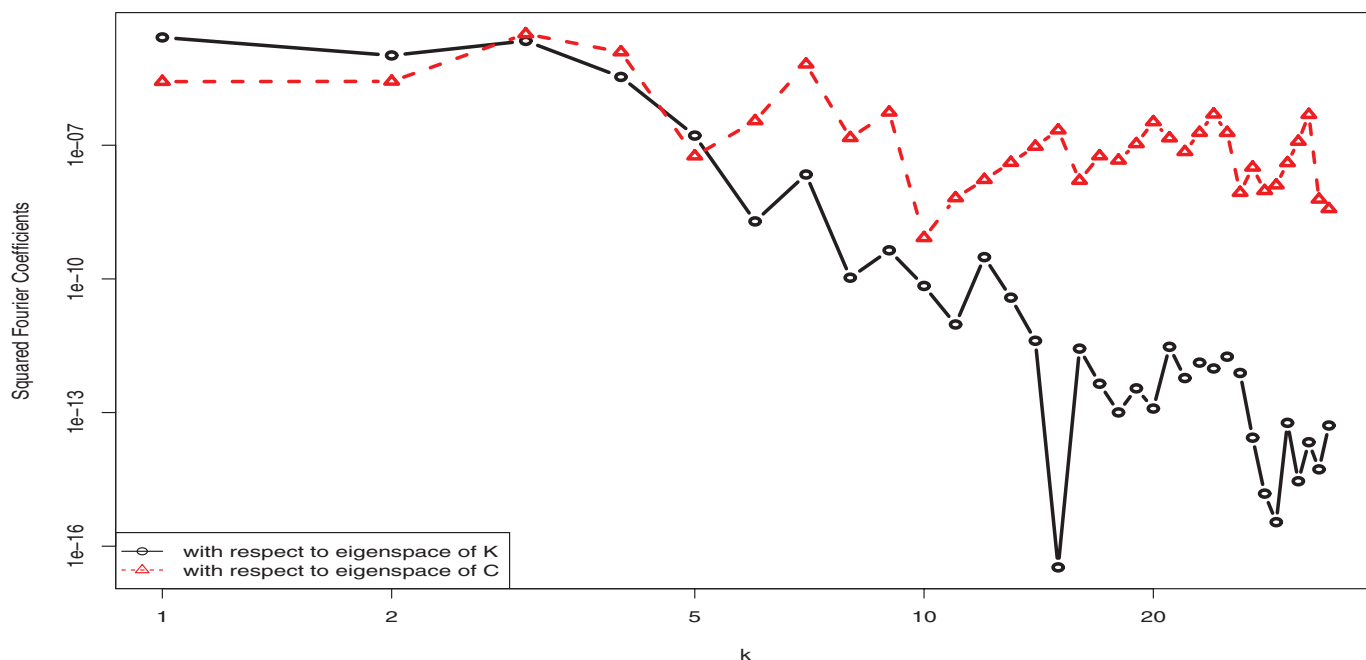


Figure 6. Canadian weather data: squared Fourier coefficients of smoothness regularization estimate of the slope function with respect to the eigenfunctions of the reproducing kernel  $K$  or  $C_n$ . Both axes are in log scale. The online version of this figure is in color.

findings are similar to those reported in Ramsay and Silverman (2005) in which a restricted basis function-based, instead of reproducing kernel-based, roughness penalization approach was used.

Note that both the functional PCA-based approach and the smoothness regularization approach assume that the slope function comes from a certain class of smooth functions where smoothness is characterized through the decay of the Fourier coefficients under a certain basis. More specifically, the functional PCA takes the Karhunen–Loève basis of  $X$ , whereas the smoothness regularization approach takes the eigenfunctions of the reproducing kernel  $K$  as the basis functions. As seen before, the assumption made by the functional PCA approach might be questionable for the current data example. To further demonstrate this and validate the appropriateness of the smoothness regularization approach, we computed the Fourier coefficients of our estimated slope function with respect to both the eigenfunctions of  $K$  and  $C_n$ . The squared coefficients are given in Figure 6. It again confirms that the assumption made by the smoothness regularization approach regarding the slope function is more reasonable.

## 5. DISCUSSION

In the present article, we establish the minimax rate of convergence for prediction with functional predictors in a general setting where the reproducing kernel  $K$  and the covariance function  $C$  are not necessarily well-aligned. The results show that the optimal prediction rate depends not only on the eigenvalues of  $K$  and  $C$ , but also on the alignment of the eigenfunctions of the two kernels. The prediction problem is more difficult if the eigenfunctions of the two kernels are not appropriately aligned. In contrast, the existing literature typically assumes that the two kernels are perfectly or nearly perfectly aligned.

We also show that the method of regularization can achieve the optimal prediction if appropriately tuned. This is to be contrasted with existing work in two ways. First, it is worth pointing out that the optimality established here for the method of regularization needs to be contrasted with the FPCA-based approaches. Because the success of these methods hinges upon the assumption that the leading principal components, although entirely determined by the predictor itself, are the most predictive of the response, it is not immediately clear to us if they can also achieve the optimal rate in general when the eigenfunctions of the reproducing kernel and covariance function differ. In addition, earlier work on the method of regularization for functional linear regression has assumed that the eigenfunctions of the reproducing kernel and covariance function are either identical or close to being identical. We show here that the optimality of this estimator extends much more generally.

To fix ideas, we have focused in this article on univariate functions, that is, the domain of the slope function and the functional predictors  $\mathcal{T} \subset \mathbb{R}$ ; and used the usual Sobolev space on  $\mathcal{T} = [0, 1]$  as a working example. Our results, however, apply to the more general RKHSs, when  $\mathcal{T}$  is a compact set in an arbitrary Euclidean space. In particular, the optimal rate of convergence and adaptivity hold true for Sobolev spaces on  $\mathcal{T} = [0, 1]^2$  with the rate of decay  $r$  determined by the corresponding reproducing kernel and covariance operator. Such a setting can have useful applications in spatial statistics and image analysis.

## 6. PROOFS

### 6.1 Proof of Theorem 1

Note that any lower bound for a specific case yields immediately a lower bound for the general case. It therefore suffices to consider the case when  $\epsilon \sim N(0, \sigma^2)$ . Denote by  $M$ , the smallest integer greater than  $c_0 n^{1/(2r+1)}$  for some constant  $c_0 > 0$  to

be specified later. For a  $\theta = (\theta_{M+1}, \dots, \theta_{2M}) \in \{0, 1\}^M$ , let

$$f_\theta = \sum_{k=M+1}^{2M} \theta_k M^{-1/2} L_{K^{1/2}} \varphi_k.$$

First, observe that  $f_\theta \in \mathcal{H}(K)$  since

$$\begin{aligned} \|f_\theta\|_{\mathcal{H}(K)}^2 &= \left\| \sum_{k=M+1}^{2M} \theta_k M^{-1/2} L_{K^{1/2}} \varphi_k \right\|_{\mathcal{H}(K)}^2 \\ &= \sum_{k=M+1}^{2M} \theta_k^2 M^{-1} \|L_{K^{1/2}} \varphi_k\|_{\mathcal{H}(K)}^2 \\ &\leq \sum_{k=M+1}^{2M} M^{-1} \|L_{K^{1/2}} \varphi_k\|_{\mathcal{H}(K)}^2 = 1, \end{aligned}$$

where we used the fact that

$$\begin{aligned} \langle L_{K^{1/2}} \varphi_j, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{H}(K)} &= \langle \varphi_j, L_K L_{K^{1/2}} \varphi_k \rangle_{\mathcal{H}(K)} \\ &= \langle \varphi_j, L_K \varphi_k \rangle_{\mathcal{H}(K)} \\ &= \langle \varphi_j, \varphi_k \rangle_{\mathcal{L}_2} = \delta_{jk}. \end{aligned}$$

The Varshamov–Gilbert bound shows that for any  $M \geq 8$ , there exists a set  $\Theta = \{\theta^{(0)}, \theta^{(1)}, \dots, \theta^{(N)}\} \subset \{0, 1\}^M$  such that

- (a)  $\theta^{(0)} = (0, \dots, 0)'$ ;
- (b)  $H(\theta, \theta') > M/8$  for any  $\theta \neq \theta' \in \Theta$ , where  $H(\cdot, \cdot)$  is the Hamming distance; and
- (c)  $N \geq 2^{M/8}$ .

We now invoke the results from Tsybakov (2009) to establish the lower bound that is based upon testing multiple hypotheses. To this end, denote by  $P_\theta$  the joint distribution of  $\{(X_i, Y_i) : i \geq 1\}$  with  $\beta_0 = f_\theta$ . First observe that for any  $\theta, \theta' \in \Theta$ ,

$$\begin{aligned} \log(P_{\theta'}/P_\theta) &= \frac{1}{\sigma^2} \sum_{i=1}^n \left( Y_i - \int X_i f_\theta \right) \int X_i (f_\theta - f_{\theta'}) \\ &\quad - \frac{1}{2\sigma^2} \sum_{i=1}^n \left[ \int X_i (f_\theta - f_{\theta'}) \right]^2. \end{aligned}$$

Therefore, the Kullback–Leibler distance between  $P_\theta$  and  $P_{\theta'}$  can be given by

$$\begin{aligned} \mathcal{K}(P_{\theta'}/P_\theta) &= \frac{n}{2\sigma^2} \|L_{C^{1/2}}(f_\theta - f_{\theta'})\|_{\mathcal{L}_2}^2 \\ &= \frac{n}{2\sigma^2} \left\| \sum_{k=M+1}^{2M} (\theta_k - \theta'_k) M^{-1/2} L_{C^{1/2}} L_{K^{1/2}} \varphi_k \right\|_{\mathcal{L}_2}^2. \end{aligned}$$

Note that

$$\begin{aligned} \langle L_{C^{1/2}} L_{K^{1/2}} \varphi_k, L_{C^{1/2}} L_{K^{1/2}} \varphi_j \rangle_{\mathcal{L}_2} &= \langle \varphi_k, L_K L_{C^{1/2}} L_{K^{1/2}} \varphi_j \rangle_{\mathcal{L}_2} \\ &= \langle \varphi_k, s_j \varphi_j \rangle_{\mathcal{L}_2} = s_j \delta_{kj}. \end{aligned}$$

Hence,

$$\begin{aligned} \mathcal{K}(P_{\theta'}/P_\theta) &= \frac{n}{2\sigma^2} \sum_{k=M+1}^{2M} M^{-1} (\theta_k - \theta'_k)^2 s_k \\ &\leq \frac{ns_M}{2M\sigma^2} \sum_{k=M+1}^{2M} (\theta_k - \theta'_k)^2 \\ &= \frac{2ns_M}{M\sigma^2} H(\theta, \theta') \end{aligned}$$

$$\begin{aligned} &\leq 2ns_M/\sigma^2 \\ &\leq 2c_2 n M^{-2r}/\sigma^2. \end{aligned}$$

This implies that for any  $0 < \alpha < 1/8$ ,

$$\frac{1}{N} \sum_{j=1}^N \mathcal{K}(P_{\theta^{(j)}} | P_{\theta^{(0)}}) \leq 2c_2 n M^{-2r}/\sigma^2 \leq \alpha \log 2^{M/8} \leq \alpha \log N \quad (11)$$

by taking  $c_0 = c\alpha^{-1/(2r+1)}$  with a large enough numerical constant  $c > 0$ .

On the other hand, for any  $\theta, \theta' \in \Theta$

$$\begin{aligned} \|L_{C^{1/2}}(f_{\theta'} - f_\theta)\|_{\mathcal{L}_2}^2 &= \left\| \sum_{k=M+1}^{2M} (\theta_k - \theta'_k) M^{-1/2} L_{C^{1/2}} L_{K^{1/2}} \varphi_k \right\|_{\mathcal{L}_2}^2 \\ &= \sum_{k=M+1}^{2M} M^{-1} (\theta_k - \theta'_k)^2 \|L_{C^{1/2}} L_{K^{1/2}} \varphi_k\|_{\mathcal{L}_2}^2 \\ &= \sum_{k=M+1}^{2M} M^{-1} (\theta_k - \theta'_k)^2 s_k \\ &\geq M^{-1} s_{2M} \sum_{k=M+1}^{2M} (\theta_k - \theta'_k)^2 \\ &= 4M^{-1} s_{2M} H(\theta, \theta') \\ &\geq s_{2M}/2 \\ &\geq c_1 2^{-(2r+1)} M^{-2r} \\ &\geq 2c\alpha^{2r/(2r+1)} n^{-\frac{2r}{2r+1}}, \end{aligned}$$

for some numerical constant  $c > 0$ . As shown by Tsybakov (2009), this, together with (11), implies that

$$\begin{aligned} \inf_{\tilde{f}} \sup_{\theta \in \Theta} P_\theta \{ \|L_{C^{1/2}}(\tilde{f} - f_\theta)\|_{\mathcal{L}_2}^2 \geq c\alpha^{2r/(2r+1)} n^{-\frac{2r}{2r+1}} \} \\ \geq \frac{\sqrt{N}}{1 + \sqrt{N}} \left( 1 - 2\alpha - \sqrt{\frac{2\alpha}{\log N}} \right). \end{aligned}$$

Therefore,

$$\begin{aligned} \liminf_{n \rightarrow \infty} \sup_{\tilde{f} \in \Theta} P_\theta \{ \|L_{C^{1/2}}(\tilde{f} - f_\theta)\|_{\mathcal{L}_2}^2 \geq c\alpha^{2r/(2r+1)} n^{-\frac{2r}{2r+1}} \} \\ \geq 1 - 2\alpha, \end{aligned}$$

which yields that

$$\lim_{a \rightarrow 0} \liminf_{n \rightarrow \infty} \sup_{\tilde{f} \in \Theta} P_\theta \{ \|L_{C^{1/2}}(\tilde{f} - f_\theta)\|_{\mathcal{L}_2}^2 \geq an^{-\frac{2r}{2r+1}} \} = 1.$$

Now the desired claim follows from the facts that  $\{f_\theta : \theta \in \Theta\} \subset \mathcal{H}(K)$ , and under  $P_\theta$ ,

$$\mathcal{E}(\tilde{\eta}) = \|L_{C^{1/2}}(\tilde{f} - f_\theta)\|_{\mathcal{L}_2}^2,$$

where

$$\tilde{\eta} = \int_T X(t) \tilde{f}(t) dt.$$

## 6.2 Proof of Theorem 2

Recall that  $L_{K^{1/2}}(\mathcal{L}_2) = \mathcal{H}(K)$ . Therefore, there exist  $f_0, \hat{f} \in \mathcal{L}_2$  such that  $\beta_0 = L_{K^{1/2}} f_0$  and  $\hat{\beta}_\lambda = L_{K^{1/2}} \hat{f}_\lambda$ . For brevity, we shall assume that  $\mathcal{H}(K)$  is dense in  $\mathcal{L}_2$ , which ensures that  $f_0$  and  $\hat{f}_\lambda$  are uniquely defined, in what follows. The proof in the

general case proceeds in exactly the same fashion by restricting ourselves to  $\mathcal{L}_2/\ker(L_{K^{1/2}})$ .

For brevity, we shall write  $T = L_{K^{1/2}}C_{K^{1/2}}$  in what follows. We shall denote by  $T^\nu$  a linear operator from  $\mathcal{L}_2$  to  $\mathcal{L}_2$  such that  $T^\nu \varphi_k = s_k^\nu \varphi_k$ . It is not hard to see that

$$\mathcal{E}(\hat{\eta}) = \|T^{1/2}(\hat{f}_\lambda - f_0)\|_{\mathcal{L}_2}^2.$$

It is also clear that

$$\hat{f}_\lambda = \operatorname{argmin}_{f \in \mathcal{L}_2} \left[ \frac{1}{n} \sum_{i=1}^n (Y_i - \langle X_i, L_{K^{1/2}} f \rangle_{\mathcal{L}_2})^2 + \lambda \|f\|_{\mathcal{L}_2}^2 \right].$$

Recall that

$$Y_i = \langle X_i, L_{K^{1/2}} f_0 \rangle_{\mathcal{L}_2} + \epsilon_i.$$

Write

$$C_n(s, t) = \frac{1}{n} \sum_{i=1}^n X_i(s)X_i(t)$$

and  $T_n = L_{K^{1/2}}L_{C_n}L_{K^{1/2}}$ , where  $L_{C_n}$  is an integral operator such that for any  $h \in \mathcal{L}_2$ ,

$$L_{C_n}h(\cdot) = \int_{\mathcal{T}} C_n(s, \cdot)h(s)ds.$$

Therefore,

$$\hat{f}_\lambda = (T_n + \lambda \mathbf{1})^{-1}(T_n f_0 + g_n)$$

where  $\mathbf{1}$  is the identity operator and

$$g_n = \frac{1}{n} \sum_{i=1}^n \epsilon_i L_{K^{1/2}} X_i.$$

Define

$$f_\lambda = (T + \lambda \mathbf{1})^{-1} T f_0.$$

By triangular inequality,

$$\|T^{1/2}(\hat{f}_\lambda - f_0)\|_{\mathcal{L}_2} = \|T^{1/2}(f_\lambda - f_0)\|_{\mathcal{L}_2} + \|T^{1/2}(\hat{f}_\lambda - f_\lambda)\|_{\mathcal{L}_2}. \tag{12}$$

The first term on the right-hand side can be easily bounded. We appeal to the following lemma. Please refer to the Appendix for proofs of all lemmas.

*Lemma 1.* For any  $0 < \nu < 1$ ,

$$\|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2} \leq (1 - \nu)^{1-\nu} \nu^\nu \lambda^\nu \|f_0\|_{\mathcal{L}_2}. \tag{13}$$

Taking  $\nu = 1/2$  in Lemma 1 gives

$$\|T^{1/2}(f_\lambda - f_0)\|_{\mathcal{L}_2}^2 \leq \frac{1}{4} \lambda \|f_0\|_{\mathcal{L}_2}^2.$$

We now turn to the second term on the right-hand side of (12). Observe that

$$\begin{aligned} f_\lambda - \hat{f}_\lambda &= (T + \lambda \mathbf{1})^{-1}(T_n + \lambda \mathbf{1})(f_\lambda - \hat{f}_\lambda) \\ &\quad + (T + \lambda \mathbf{1})^{-1}(T - T_n)(f_\lambda - \hat{f}_\lambda). \end{aligned}$$

Recall that

$$(T_n + \lambda \mathbf{1})\hat{f}_\lambda = T_n f_0 - g_n.$$

Therefore,

$$\begin{aligned} f_\lambda - \hat{f}_\lambda &= (T + \lambda \mathbf{1})^{-1} T_n (f_\lambda - f_0) + \lambda (T + \lambda \mathbf{1})^{-1} f_\lambda \\ &\quad + (T + \lambda \mathbf{1})^{-1} g_n + (T + \lambda \mathbf{1})^{-1} (T - T_n)(f_\lambda - \hat{f}_\lambda) \\ &= (T + \lambda \mathbf{1})^{-1} T_n (f_\lambda - f_0) + \lambda T f_0 + (T + \lambda \mathbf{1})^{-1} g_n \\ &\quad + (T + \lambda \mathbf{1})^{-1} (T - T_n)(f_\lambda - \hat{f}_\lambda) \\ &= (T + \lambda \mathbf{1})^{-1} T (f_\lambda - f_0) + (T + \lambda \mathbf{1})^{-1} (T_n - T) \\ &\quad \times (f_\lambda - f_0) + \lambda T f_0 + (T + \lambda \mathbf{1})^{-1} g_n \\ &\quad + (T + \lambda \mathbf{1})^{-1} (T - T_n)(f_\lambda - \hat{f}_\lambda). \end{aligned}$$

We first consider bounding  $\|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2}$  for some  $0 < \nu < 1/2 - 1/(4r)$ . By triangular inequality,

$$\begin{aligned} \|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} &\leq \|T^\nu(T + \lambda \mathbf{1})^{-1} T (f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\quad + \|T^\nu(T + \lambda \mathbf{1})^{-1} (T_n - T)(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\quad + \lambda \|T^{1+\nu} f_0\|_{\mathcal{L}_2} + \|T^\nu(T + \lambda \mathbf{1})^{-1} g_n\|_{\mathcal{L}_2} \\ &\quad + \|T^\nu(T + \lambda \mathbf{1})^{-1} (T - T_n)(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2}. \end{aligned}$$

Next, we make use of another auxiliary lemma.

*Lemma 2.* Assume that there exists a constant  $c_3 > 0$  such that for any  $f \in \mathcal{L}_2$

$$\mathbb{E}\langle X, f \rangle_{\mathcal{L}_2}^4 \leq c_3 (\mathbb{E}\langle X, f \rangle_{\mathcal{L}_2}^2)^2. \tag{14}$$

Then, for any  $\nu > 0$  such that  $2r(1 - 2\nu) > 1$ ,

$$\|T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} = O_p\left(\left(n\lambda^{1-2\nu+1/(2r)}\right)^{-1/2}\right),$$

where  $\|\cdot\|_{\text{op}}$  stands for the usual operator norm, that is,  $\|U\|_{\text{op}} = \sup_{h: \|h\|_{\mathcal{L}_2}=1} \|Uh\|_{\mathcal{L}_2}$  for an operator  $U: \mathcal{L}_2 \mapsto \mathcal{L}_2$ .

An application of Lemma 2 yields

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1}(T - T_n)(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} &\leq \|T^\nu(T + \lambda \mathbf{1})^{-1}(T - T_n)T^{-\nu}\|_{\text{op}} \|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} \\ &\leq o_p(1) \|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} \end{aligned}$$

whenever  $\lambda \geq cn^{-2r/(2r+1)}$  for some constant  $c > 0$ . Similarly,

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)(f_\lambda - f_0)\|_{\mathcal{L}_2} &\leq \|T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} \|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\leq o_p(1) \|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2}. \end{aligned}$$

Therefore,

$$\begin{aligned} \|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} &= O_p\left(\|T^\nu(T + \lambda \mathbf{1})^{-1} T (f_\lambda - f_0)\|_{\mathcal{L}_2} \right. \\ &\quad \left. + \lambda \|T^{1+\nu} f_0\| + \|T^\nu(T + \lambda \mathbf{1})^{-1} g_n\|_{\mathcal{L}_2}\right). \end{aligned}$$

By Lemma 1,

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1} T (f_\lambda - f_0)\|_{\mathcal{L}_2} &\leq \|T^\nu(T + \lambda \mathbf{1})^{-1} T^{1-\nu}\|_{\text{op}} \|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\leq \|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\leq (1 - \nu)^{1-\nu} \nu^\nu \lambda^\nu \|f_0\|_{\mathcal{L}_2}. \end{aligned}$$

Together with Lemma 3 stated later, we conclude that

$$\|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} = O_p\left(\lambda^\nu + (n\lambda^{1-2\nu+1/(2r)})^{-1/2}\right) = O_p(\lambda^\nu)$$

provided that  $c_1 n^{-2r/(2r+1)} \leq \lambda \leq c_2 n^{-2r/(2r+1)}$  for some constants  $0 < c_1 < c_2 < \infty$ .

*Lemma 3.* For any  $0 \leq \nu \leq 1/2$ ,

$$\|T^\nu(T + \lambda \mathbf{1})^{-1} g_n\|_{\mathcal{L}_2} = O_p\left(\left(n\lambda^{1-2\nu+1/(2r)}\right)^{-1/2}\right).$$

We are now in position to bound  $\|T^{1/2}(f_\lambda - \hat{f}_\lambda)\|$ . Recall that

$$\begin{aligned} \|T^{1/2}(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} &\leq \|T^{1/2}(T + \lambda\mathbf{1})^{-1}T(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\quad + \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\quad + \lambda\|T^{1+1/2}f_0\| + \|T^{1/2}(T + \lambda\mathbf{1})^{-1}g_n\|_{\mathcal{L}_2} \\ &\quad + \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T - T_n)(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2}. \end{aligned}$$

We now bound the five terms on the right-hand side separately. By Lemma 1,

$$\begin{aligned} \|T^{1/2}(T + \lambda\mathbf{1})^{-1}T(f_\lambda - f_0)\|_{\mathcal{L}_2} &\leq \|T^{1/2}(T + \lambda\mathbf{1})^{-1}T^{1/2}\|_{\text{op}} \\ &\quad \times \|T^{1/2}(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\leq \frac{1}{2}\lambda^{1/2}\|f_0\|_{\mathcal{L}_2}. \end{aligned}$$

We appeal to the following result.

*Lemma 4.* Under the conditions of Lemma 2,

$$\|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} = O_p\left((n\lambda^{1/(2r)})^{-1/2}\right).$$

By Lemmas 1 and 4,

$$\begin{aligned} \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)(f_\lambda - f_0)\|_{\mathcal{L}_2} &\leq \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}}\|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2} \\ &\leq O_p\left((n\lambda^{1/(2r)})^{-1/2}\lambda^\nu\right) = o_p\left((n\lambda^{1/(2r)})^{-1/2}\right). \end{aligned}$$

Similarly,

$$\begin{aligned} \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} &\leq \|T^{1/2}(T + \lambda\mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}}\|T^\nu(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} \\ &\leq O_p\left((n\lambda^{1/(2r)})^{-1/2}\lambda^\nu\right) = o_p\left((n\lambda^{1/(2r)})^{-1/2}\right). \end{aligned}$$

By Lemma 3,

$$\|T^{1/2}(T + \lambda\mathbf{1})^{-1}g_n\|_{\mathcal{L}_2} = O_p\left((n\lambda^{1/(2r)})^{-1/2}\right).$$

Together with the fact that  $\lambda\|T^{1+1/2}f_0\| = O(\lambda)$ , we conclude that

$$\|T^{1/2}(f_\lambda - \hat{f}_\lambda)\|_{\mathcal{L}_2} = O_p\left(n^{-\frac{2r}{2r+1}}\right).$$

### 6.3 Proof of Theorem 3

Consider the following Rademacher type of process:

$$R_n(b) = \frac{1}{n} \sum_{i=1}^n w_i \langle X_i, b \rangle_{\mathcal{L}_2},$$

where  $w_i$ 's are iid Rademacher random variables, that is,  $\mathbb{P}(w_i = 1) = \mathbb{P}(w_i = -1) = 1/2$ . Define

$$\|R_n\|_{\mathcal{B}(\delta)} = \sup_{b \in \mathcal{B}(\delta)} |R_n(b)|,$$

where

$$\mathcal{B}(\delta) = \{b \in \mathcal{H}(K) : \|b\|_{\mathcal{H}(K)} \leq 1 \text{ and } \|L_{C^{1/2}}b\|_{\mathcal{L}_2} \leq \delta\}.$$

Define

$$\rho_1 = \inf\{\rho \geq n^{-1} \log n : \mathbb{E}\|R_n\|_{\mathcal{B}(\delta)} \leq \rho^{1/2}\delta + \rho, \forall \delta \in [0, 1]\}.$$

Similarly, write

$$\hat{\rho}_1 = \inf\{\rho \geq n^{-1} \log n : \mathbb{E}_w\|R_n\|_{\mathcal{B}(\delta)} \leq \rho^{1/2}\delta + \rho, \forall \delta \in [0, 1]\},$$

where  $\mathbb{E}_w$  stands for expectation taken over the Rademacher random variables  $w_i$ 's only. We first note the following result.

*Lemma 5.* Under the condition of Theorem 3, there exist constants  $c_1, c_2, c_3 > 0$  such that

$$c_1\gamma_n(\delta) - c_2n^{-1}(\log n) \leq \mathbb{E}\|R_n\|_{\mathcal{B}} \leq c_3\gamma_n(\delta).$$

It is clear from Lemma 5 that  $0 < \inf \rho_1/\rho \leq \sup \rho_1/\rho < \infty$ . Following the same argument, it can be shown that  $0 < \inf \hat{\rho}_1/\hat{\rho} \leq \sup \hat{\rho}_1/\hat{\rho} < \infty$ . It now suffices to show that  $\hat{\rho}_1/\rho_1$  is also bounded away from 0 and  $+\infty$ .

For  $b \in \mathcal{B}(\delta)$ , let  $f = L_{K^{-1/2}}b$ . Then

$$\begin{aligned} |R_n(b)| &= \left| \frac{1}{n} \sum_{i=1}^n w_i \langle X_i, b \rangle_{\mathcal{L}_2} \right| \\ &= \left| \frac{1}{n} \sum_{i=1}^n w_i \langle L_{K^{1/2}}X_i, f \rangle_{\mathcal{L}_2} \right| \\ &\leq \frac{1}{n} \sum_{i=1}^n |\langle L_{K^{1/2}}X_i, f \rangle_{\mathcal{L}_2}|. \end{aligned}$$

By Cauchy–Schwartz inequality, this can be further bounded.

$$|R_n(b)| \leq \frac{1}{n} \sum_{i=1}^n \|L_{K^{1/2}}X_i\|_{\mathcal{L}_2} \|f\|_{\mathcal{L}_2} \leq \frac{1}{n} \sum_{i=1}^n \|L_{K^{1/2}}X_i\|_{\mathcal{L}_2},$$

where the second inequality follows from the fact that  $\|f\|_{\mathcal{L}_2} = \|b\|_{\mathcal{H}(K)} \leq 1$ . Note that the rightmost hand side converges almost surely to  $\mathbb{E}\|L_{K^{1/2}}X\|_{\mathcal{L}_2} < \infty$  by strong law of large numbers. On the other hand,

$$\mathbb{E}\|R_n(b)\|^2 = \frac{1}{n} \mathbb{E}(w_i \langle X_i, b \rangle_{\mathcal{L}_2})^2 = \|L_{C^{1/2}}b\|_{\mathcal{L}_2}^2 \leq \delta^2.$$

The rest of the proof follows in the same fashion as that of Koltchinskii and Yuan (2010, sec. 3.2) and is therefore omitted.

## APPENDIX: AUXILIARY RESULTS

### Proof of Lemma 1

Write

$$f_0 = \sum_{k \geq 1} a_k \varphi_k.$$

Then,

$$f_\lambda = \sum_{k \geq 1} \frac{s_k a_k}{\lambda + s_k} \varphi_k.$$

Therefore,

$$\|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2}^2 = \sum_{k \geq 1} s_k^{2\nu} \left( \frac{\lambda a_k}{\lambda + s_k} \right)^2 \leq \max_{k \geq 1} \frac{\lambda^2 s_k^{2\nu}}{(\lambda + s_k)^2} \sum_{k \geq 1} a_k^2.$$

By Young's inequality,  $\lambda + s_k \geq (1-\nu)^{-1}\nu^{-\nu}\lambda^{1-\nu}s_k^\nu$ . Hence,

$$\|T^\nu(f_\lambda - f_0)\|_{\mathcal{L}_2}^2 \leq (1-\nu)^{2(1-\nu)}\nu^{2\nu}\lambda^{2\nu}\|f_0\|_{\mathcal{L}_2}^2.$$

### Proof of Lemma 2

Recall that

$$\begin{aligned} \|T^\nu(T + \lambda\mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} &= \sup_{h: \|h\|_{\mathcal{L}_2}=1} |\langle h, T^\nu(T + \lambda\mathbf{1})^{-1}(T_n - T)T^{-\nu}h \rangle_{\mathcal{L}_2}|. \end{aligned}$$

Write

$$h = \sum_{k \geq 1} h_k \varphi_k.$$

Then,

$$\begin{aligned} \langle h, T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}h \rangle_{\mathcal{L}_2} &= \langle T^\nu(T + \lambda \mathbf{1})^{-1}h, (T_n - T)T^{-\nu}h \rangle_{\mathcal{L}_2} \\ &= \left\langle \sum_{j \geq 1} \frac{s_j^\nu h_j}{s_j + \lambda} \varphi_j, \sum_{k \geq 1} s_k^{-\nu} h_k (T_n - T) \varphi_k \right\rangle_{\mathcal{L}_2} \\ &= \sum_{j,k \geq 1} \frac{s_j^\nu s_k^{-\nu} h_j h_k}{s_j + \lambda} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}. \end{aligned}$$

An application of Cauchy–Schwartz inequality yields

$$\begin{aligned} &\left| \sum_{j,k \geq 1} \frac{s_j^\nu s_k^{-\nu} h_j h_k}{s_j + \lambda} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2} \right| \\ &\leq \left( \sum_{j,k \geq 1} h_j^2 h_k^2 \right)^{1/2} \left( \sum_{j,k \geq 1} \frac{s_j^{2\nu} s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2}. \end{aligned}$$

Hence,

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} &\leq \left( \sum_{j,k \geq 1} \frac{s_j^{2\nu} s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2}. \quad (15) \end{aligned}$$

Now consider the expectation of the right-hand side. By Jensen’s inequality,

$$\begin{aligned} \mathbb{E} \left( \sum_{j,k \geq 1} \frac{s_j^{2\nu} s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2} \\ \leq \left( \sum_{j,k \geq 1} \frac{s_j^{2\nu} s_k^{-2\nu}}{(s_j + \lambda)^2} \mathbb{E} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2}. \end{aligned}$$

Note that

$$\begin{aligned} \mathbb{E} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 &= \mathbb{E} \langle L_{K^{1/2}} \varphi_j, (L_{C_n} - L_C) L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2}^2 \\ &= \mathbb{E} \left( \int_{\mathcal{T}^2} (L_{K^{1/2}} \varphi_j)(s) (C_n(s, t) - C(s, t)) (L_{K^{1/2}} \varphi_k)(t) \right)^2 \\ &= \mathbb{E} \left( \frac{1}{n} \sum_{i=1}^n \int_{\mathcal{T}^2} (L_{K^{1/2}} \varphi_j)(s) (X_i(s) X_i(t) - \mathbb{E} X_i(s) X_i(t)) (L_{K^{1/2}} \varphi_k)(t) \right)^2 \\ &= \frac{1}{n} \mathbb{E} \left( \int_{\mathcal{T}^2} (L_{K^{1/2}} \varphi_j)(s) (X(s) X(t) - \mathbb{E} X(s) X(t)) (L_{K^{1/2}} \varphi_k)(t) \right)^2 \\ &\leq \frac{1}{n} \mathbb{E} \left( \int_{\mathcal{T}^2} (L_{K^{1/2}} \varphi_j)(s) X(s) X(t) (L_{K^{1/2}} \varphi_k)(t) \right)^2. \end{aligned}$$

An application of Cauchy–Schwartz inequality yields

$$\begin{aligned} \mathbb{E} \langle L_{K^{1/2}} \varphi_j, (L_{C_n} - L_C) L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2}^2 &\leq \frac{1}{n} \mathbb{E}^{1/2} \left( \int_{\mathcal{T}} (L_{K^{1/2}} \varphi_j)(t) X(t) dt \right)^4 \mathbb{E}^{1/2} \left( \int_{\mathcal{T}} (L_{K^{1/2}} \varphi_k)(t) X(t) dt \right)^4 \\ &\leq c_3 n^{-1} \mathbb{E} \left( \int_{\mathcal{T}} (L_{K^{1/2}} \varphi_j)(t) X(t) dt \right)^2 \mathbb{E} \left( \int_{\mathcal{T}} (L_{K^{1/2}} \varphi_k)(t) X(t) dt \right)^2 \\ &= c_3 n^{-1} \|T^{1/2} \varphi_j\|_{\mathcal{L}_2}^2 \|T^{1/2} \varphi_k\|_{\mathcal{L}_2}^2 = c_3 n^{-1} s_j s_k. \end{aligned}$$

Therefore,

$$\mathbb{E} \left( \sum_{j,k \geq 1} \frac{s_j^{2\nu} s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T) \varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2} \leq \left( \frac{1}{n} \sum_{j,k \geq 1} \frac{s_j^{1+2\nu} s_k^{1-2\nu}}{(s_j + \lambda)^2} \right)^{1/2}. \quad (16)$$

Note that  $s_k^{1-2\nu}$  is summable because  $(1 - 2\nu)(2r) > 1$ . We now appeal to the following lemma.

*Lemma 6.* If there exist constants  $0 < c_1 < c_2 < \infty$  such that  $c_1 k^{-2r} < s_k < c_2 k^{-2r}$ , then there exist constants  $c_3, c_4 > 0$  depending only on  $c_1, c_2$  such that

$$c_4 \lambda^{-1/(2r)} \leq \sum_{j \geq 1} \frac{s_j^{1+2\nu}}{(\lambda + s_j)^{1+2\nu}} \leq c_3 (1 + \lambda^{-1/(2r)}).$$

Thus, by Lemma 6,

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} &\leq c (n \lambda^{1-2\nu})^{-1/2} \left( \sum_{j \geq 1} \frac{s_j^{1+2\nu}}{(\lambda + s_j)^{1+2\nu}} \right)^{1/2} \\ &\leq c (n \lambda^{1-2\nu+1/(2r)})^{-1/2}. \end{aligned}$$

The proof is now completed by Markov inequality.

**Proof of Lemma 3**

Recall that

$$g_n = \frac{1}{n} \sum_{i=1}^n \epsilon_i L_{K^{1/2}} X_i.$$

Therefore,

$$\begin{aligned} \|T^\nu(T + \lambda \mathbf{1})^{-1} g_n\|_{\mathcal{L}_2}^2 &= \sum_{k \geq 1} \langle T^\nu(T + \lambda \mathbf{1})^{-1} g_n, \varphi_k \rangle_{\mathcal{L}_2}^2 \\ &= \sum_{k \geq 1} \langle g_n, (T + \lambda \mathbf{1})^{-1} T^\nu \varphi_k \rangle_{\mathcal{L}_2}^2 \\ &= \sum_{k \geq 1} \left\langle \frac{1}{n} \sum_{i=1}^n \epsilon_i L_{K^{1/2}} X_i, \frac{s_k^\nu}{\lambda + s_k} \varphi_k \right\rangle_{\mathcal{L}_2}^2 \\ &= \sum_{k \geq 1} \frac{s_k^{2\nu}}{(\lambda + s_k)^2} \left( \frac{1}{n} \sum_{i=1}^n \epsilon_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2} \right)^2. \end{aligned}$$

Note that

$$\mathbb{E} (\epsilon_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2}) = \mathbb{E} (\epsilon_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2} | X_i) = 0.$$

Hence,

$$\begin{aligned} \mathbb{E} \|T^\nu(T + \lambda \mathbf{1})^{-1} g_n\|_{\mathcal{L}_2}^2 &= \sum_{k \geq 1} \frac{s_k^{2\nu}}{(\lambda + s_k)^2} \mathbb{E} \left( \frac{1}{n} \sum_{i=1}^n \epsilon_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2} \right)^2 \\ &= \frac{1}{n} \sum_{k \geq 1} \frac{s_k^{2\nu}}{(\lambda + s_k)^2} \mathbb{E} (\epsilon_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2})^2 \\ &= \frac{\sigma^2}{n} \sum_{k \geq 1} \frac{s_k^{2\nu}}{(\lambda + s_k)^2} \mathbb{E} \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2}^2 \\ &= \frac{\sigma^2}{n} \sum_{k \geq 1} \frac{s_k^{1+2\nu}}{(\lambda + s_k)^2} \\ &\leq \frac{\sigma^2}{n \lambda^{1-2\nu}} \sum_{k \geq 1} \frac{s_k^{1+2\nu}}{(\lambda + s_k)^{1+2\nu}} \\ &\leq \frac{c \sigma^2}{n \lambda^{1-2\nu+1/(2r)}}. \end{aligned}$$

The proof can now be completed by Markov inequality.

**Proof of Lemma 4**

Similar to (15), by Cauchy–Schwartz inequality,

$$\begin{aligned} & \|T^{1/2}(T + \lambda \mathbf{1})^{-1}(T_n - T)T^{-\nu}\|_{\text{op}} \\ &= \sup_{h: \|h\|_{\mathcal{L}_2}=1} \left| \sum_{j,k \geq 1} \frac{s_j s_k^{-\nu} h_j h_k}{s_j + \lambda} \langle \varphi_j, (T_n - T)\varphi_k \rangle_{\mathcal{L}_2} \right| \\ &\leq \left( \sum_{j,k \geq 1} \frac{s_j s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T)\varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2}. \end{aligned}$$

Following a similar argument as that of (16),

$$\mathbb{E} \left( \sum_{j,k \geq 1} \frac{s_j s_k^{-2\nu}}{(s_j + \lambda)^2} \langle \varphi_j, (T_n - T)\varphi_k \rangle_{\mathcal{L}_2}^2 \right)^{1/2} \leq \left( \frac{1}{n} \sum_{j,k \geq 1} \frac{s_j^2 s_k^{1-2\nu}}{(s_j + \lambda)^2} \right)^{1/2} \leq c(n\lambda^{1/(2\nu)})^{-1/2}.$$

The proof is now completed by Markov inequality.

**Proof of Lemma 5**

It is clear that  $\mathcal{B}(\delta) = L_{K^{1/2}}(\mathcal{F}(\delta))$ , where

$$\mathcal{F}(\delta) = \{f \in \mathcal{L}_2 : \|f\|_{\mathcal{L}_2} \leq 1 \text{ and } \|T^{1/2} f\|_{\mathcal{L}_2}^2 \leq \delta^2\}.$$

Denote

$$\mathcal{G} = \left\{ \sum_{k \geq 1} f_k \varphi_k : \sum_{k \geq 1} \left( \frac{f_k}{\min\{1, \delta/\sqrt{s_k}\}} \right)^2 \leq 1 \right\}.$$

It can be easily checked that  $\mathcal{G} \subset \mathcal{F} \subset \sqrt{2}\mathcal{G}$ . Therefore,

$$\sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)| \leq \|R_n\|_B \leq \sqrt{2} \sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)|.$$

By Jensen’s inequality,

$$\mathbb{E} \sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)| \leq \left( \mathbb{E} \sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)|^2 \right)^{1/2}.$$

By Cauchy–Schwartz inequality,

$$\begin{aligned} & |R_n(L_{K^{1/2}} f)|^2 \\ &= \left| \sum_{k \geq 1} f_k R_n(L_{K^{1/2}} \varphi_k) \right|^2 \\ &\leq \left( \sum_{k \geq 1} \frac{f_k^2}{\min\{1, \delta^2/s_k\}} \right) \left( \sum_{k \geq 1} \min\{1, \delta^2/s_k\} R_n^2(L_{K^{1/2}} \varphi_k) \right). \end{aligned}$$

Therefore,

$$\sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)|^2 \leq \left( \sum_{k \geq 1} \min\{1, \delta^2/s_k\} R_n^2(L_{K^{1/2}} \varphi_k) \right).$$

Observe that

$$\begin{aligned} \mathbb{E} R_n^2(L_{K^{1/2}} \varphi_k) &= \mathbb{E} \left( \frac{1}{n} \sum_{i=1}^n w_i \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2} \right)^2 \\ &= \frac{1}{n} \mathbb{E} \langle X_i, L_{K^{1/2}} \varphi_k \rangle_{\mathcal{L}_2}^2 = n^{-1} s_k. \end{aligned}$$

Thus,

$$\mathbb{E} \left( \sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)|^2 \right) \leq \left( \sum_{k \geq 1} \min\{1, \delta^2/s_k\} \mathbb{E} R_n^2(L_{K^{1/2}} \varphi_k) \right) = \gamma_n^2(\delta),$$

which implies that

$$\mathbb{E} \|R_n\|_B \leq \sqrt{2} \gamma_n(\delta).$$

To prove the lower bound, we appeal to Hoffman–Jørgensen inequality that suggests that

$$\mathbb{E}^{1/2} \|R_n\|_B^2 \leq c \left( \mathbb{E} \|R_n\|_B + \frac{1}{n} \mathbb{E}^{1/2} \max_{1 \leq i \leq n} \sup_{\beta \in \mathcal{F}} \langle X_i, \beta \rangle_{\mathcal{L}_2}^2 \right).$$

Observe that

$$\begin{aligned} \mathbb{E}^{1/2} \max_{1 \leq i \leq n} \sup_{\beta \in \mathcal{F}} \langle X_i, \beta \rangle_{\mathcal{L}_2}^2 &= \mathbb{E}^{1/2} \max_{1 \leq i \leq n} \sup_{f \in \mathcal{G}} \langle L_{K^{1/2}} X_i, f \rangle_{\mathcal{L}_2}^2 \\ &\leq \mathbb{E}^{1/2} \max_{1 \leq i \leq n} \|L_{K^{1/2}} X_i\|_{\mathcal{L}_2}^2. \end{aligned}$$

Because  $\|L_{K^{1/2}} X\|_{\mathcal{L}_2}$  has exponential tails, it can be further bounded by  $c \log n$  for some constant  $c > 0$ .

Hence,

$$\begin{aligned} \mathbb{E} \|R_n\|_B &\geq c_1 \mathbb{E}^{1/2} \|R_n\|_B^2 - c_2 n^{-1} (\log n) \\ &\geq c_1 \mathbb{E}^{1/2} \left( \sup_{f \in \mathcal{G}} |R_n(L_{K^{1/2}} f)|^2 \right) - c_2 n^{-1} (\log n) \\ &\geq c_1 \gamma_n(\delta) - c_2 n^{-1} (\log n). \end{aligned}$$

**Proof of Lemma 6**

Note that

$$\begin{aligned} \sum_{k \geq 1} \frac{s_k^{1+2\nu}}{(\lambda + s_k)^{1+2\nu}} &\leq \sum_{k \geq 1} \frac{(c_1 k^{-2r})^{1+2\nu}}{(\lambda + c_2 k^{-2r})^{1+2\nu}} \\ &= c_1^{-2r(1+2\nu)} \sum_{k \geq 1} \frac{1}{(c_2 + \lambda k^{2r})^{1+2\nu}} \\ &\leq c_1^{-2r(1+2\nu)} \left( \frac{1}{c_2} + \int_1^\infty \frac{dx}{(c_2 + \lambda x^{2r})^{1+2\nu}} \right) \\ &= c_1^{-2r(1+2\nu)} \left( \frac{1}{c_2} + \lambda^{-\frac{1}{2r}} \int_{\lambda^{1/(2r)}}^\infty \frac{dy}{(c_2 + y^{2r})^{1+2\nu}} \right) \\ &\leq c_3 (1 + \lambda^{-1/(2r)}). \end{aligned}$$

Similarly,

$$\begin{aligned} \sum_{k \geq 1} \frac{s_k^{1+2\nu}}{(\lambda + s_k)^{1+2\nu}} &\geq c_2^{-2r(1+2\nu)} \sum_{k \geq 1} \frac{1}{(c_1 + \lambda k^{2r})^{1+2\nu}} \\ &\geq c_2^{-2r(1+2\nu)} \int_1^\infty \frac{dx}{(c_1 + \lambda x^{2r})^{1+2\nu}} \\ &= c_2^{-2r(1+2\nu)} \lambda^{-\frac{1}{2r}} \int_{\lambda^{1/(2r)}}^\infty \frac{dy}{(c_1 + y^{2r})^{1+2\nu}} \\ &\geq c_4 \lambda^{-1/(2r)}. \end{aligned}$$

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