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# Prediction in function-on-function linear model with partially observed functional covariate and response

Christophe Crambes<sup>†</sup> and Chayma Daayeb<sup>†,‡</sup> and Ali Gannoun<sup>†</sup> and Yousri Henchiri<sup>‡,§,\*</sup>

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## Abstract

In this work, we are interested in a function-on-function linear model in which the response and the covariate are partially observed curves. First, we reconstruct the missing part of the covariate using the observed parts. Then, we consider two strategies for dealing with the missing part of the response. The first one consists in a reconstruction in the same way as for the covariate. The second one uses regression imputation. Once the dataset is reconstructed, we estimate the slope function and give the mean square prediction error for a new observation of the covariate. Both methods are compared from a theoretical and a practical point of view.

**Keywords.** Functional linear model, Missing parts, Imputation, Partially observed functional data, Functional Principal Components, Reconstruction operator.

## 1 Introduction

Functional data analysis (FDA) is becoming progressively more and more important in Statistic (Bosq, 2000; Ramsay and Silverman, 2005; Ferraty and Vieu, 2006; Hsing and Eubank, 2015; Kokoszka and Reimherr, 2018). The functional linear model is specially a very popular model in both theoretical and applied research such as climatology, meteorology, economy, image analysis and many other fields. There is a large amount of work done on the functional linear model. In the case where the response is a real variable, the model has been widely studied (e.g. Cardot et al., 2003; Cai and Hall, 2006; Li and Hsing, 2007; Hall and Horowitz, 2007; Crambes et al., 2009; Comte and Johannes, 2012; Cai and Yuan, 2012; Brunel et al.,

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2016). In contrast, fewer researchers have tackled the problem of function-on-function linear models where the covariate and the response are both functional (e.g. Ramsay and Silverman, 2005; Yao et al., 2005; Prchal and Sarda., 2007; Aguilera et al., 2008; Lian, 2011; Ferraty et al., 2012; Crambes and Mas, 2013; Crambes et al., 2016; Luo and Qi, 2017; Benatia et al., 2017; Imaizumi and Kato, 2018; Sun et al., 2018).

In recent years, applications producing partially observed functional data have emerged. Sometimes each individual trajectory is collected only over individual specific subintervals, densely or sparsely, within the whole domain of interest. Several recent works have begun addressing the estimation of covariance functions for short functional segments observed at sparse and irregular grid points, called *functional snippets* (Lin and Wang, 2022; Lin et al., 2021) or for *fragmented functional data* observed on small subintervals (Delaigle et al., 2020). For densely observed partial data, existing studies have focused on estimating the unobserved part of curves (Kneip and Liebl, 2020; Kraus and Stefanucci, 2020), prediction (Goldberg et al., 2014), classification (Kraus and Stefanucci, 2018; Park and Simpson, 2019), functional regression (Gellar et al., 2014), and inferences (Kraus, 2019; Park et al., 2022).

To our knowledge, few articles investigate the theoretical properties of the slope operator or kernel estimators in the framework of partially observed data. In the framework of a real response, we can notice the works from Crambes and HENCHIRI (2019) and Crambes et al. (2023). The objective of this paper is to study the prediction problem when the covariate and the response are partially observed in the function-on-function linear regression setting, for which we propose two methods. The paper is organized as follows. We present the model with fully observed functional data in section 2. In section 3, we study the case of partially observed covariate and response. We reconstruct the missing parts of the functional explanatory variable and the functional response using the work Kneip and Liebl (2020), and we get theoretical results for the prediction error rate. In section 4, we propose an alternative method which consists in imputing the functional response after having reconstructed the missing parts of the functional covariate as in the previous section. We also obtain a convergence rate for the mean square prediction error. In section 5, we conduct a numerical study over simulated data in order to compare the methods in practice. Section 6 is devoted to a real dataset application. Finally, all the proofs are postponed to section 7.

## 2 The centered function-on-function linear model

### 2.1 Functional principal components regression

Let  $f : \mathcal{T} \rightarrow \mathbb{R}$  and  $g : \mathcal{T} \times \mathcal{S} \rightarrow \mathbb{R}$  be two square integrable functions defined in the Hilbert space  $\mathbb{L}^2(\mathcal{T})$  (resp.  $\mathbb{L}^2(\mathcal{T} \times \mathcal{S})$ ) i.e. the space of square integrable functions on the interval  $\mathcal{T} \subseteq \mathbb{R}$  (resp.  $\mathcal{S} \subseteq \mathbb{R}$ ). For all  $t \in \mathcal{T}$  (resp.  $s \in \mathcal{S}$ ), we set  $\|f\| = \left( \int_{\mathcal{T}} f^2(t) dt \right)^{1/2}$  and  $\|g\| = \left( \int_{\mathcal{T}} \int_{\mathcal{S}} g^2(t, s) dt ds \right)^{1/2}$ .

We use the following notation for the tensor product, defining  $u \otimes v : \mathbb{L}^2(\mathcal{T}) \rightarrow \mathbb{L}^2(\mathcal{T})$ , by  $u \otimes v = \langle u, \cdot \rangle v$ , for any functions  $u, v \in \mathbb{L}^2(\mathcal{T})$ . We define the Hilbert-Schmidt integral operator  $G : \mathbb{L}^2(\mathcal{T} \times \mathcal{S}) \rightarrow \mathbb{L}^2(\mathcal{T})$  by  $(G \cdot u)(s) = \int_{\mathcal{T}} u(t)g(t, s)dt$  for all  $g$  belonging to  $\mathbb{L}^2(\mathcal{T} \times \mathcal{S})$  and  $\int_{\mathcal{T}} \int_{\mathcal{S}} |g(t, s)|^2 dt ds < \infty$ . The function  $g$  is the Hilbert-Schmidt kernel corresponding to the operator  $G$ . Moreover, the operator  $G$  is continuous (i.e bounded) and compact.

We assume here that  $\mathcal{X}$  takes values in the space  $\mathbb{L}^2(\mathcal{T})$  and  $\mathcal{Y}$  in  $\mathbb{L}^2(\mathcal{S})$ ,  $\mathcal{X} \triangleq \{\mathcal{X}(t) \mid t \in \mathcal{T}\}$  is the predictor variable and  $\mathcal{Y} \triangleq \{\mathcal{Y}(s) \mid s \in \mathcal{S}\}$  the response variable. We observe a sample  $\{(\mathcal{X}_i, \mathcal{Y}_i)\}_{i=1}^n$ , of identically distributed and independent copies  $(\mathcal{X}, \mathcal{Y})$ , with  $\mathbb{E}(\|\mathcal{X}\|^2) < \infty$  and  $\mathbb{E}(\|\mathcal{Y}\|^2) < \infty$ . In practice, functional responses  $\mathcal{Y}_i(s)$  and functional covariates  $\mathcal{X}_i(t)$  are observed on grid points  $\mathbf{s}_i = (s_{i1}, \dots, s_{iq_i})$  and  $\mathbf{t}_i = (t_{i1}, \dots, t_{ip_i})$ . For the sake of simplicity, we assume those to be identical vectors  $\mathbf{s}, \mathbf{t}$  with lengths  $q$  and  $p$ , respectively, for each observation  $i$ .

We consider a centered function-on-function linear regression model (FFLRM) characterizing a linear relationship between the functional response and the functional predictor

$$[\mathcal{Y} - \mathbb{E}[\mathcal{Y}]](s) = \int_{\mathcal{T}} [\mathcal{X} - \mathbb{E}[\mathcal{X}]](t)\theta(t, s)dt + \epsilon(s),$$

where the bivariate functional coefficient  $(t, s) \mapsto \theta(t, s)$  is assumed to be in  $\mathbb{L}^2(\mathcal{T} \times \mathcal{S})$ , that is,  $\int_{\mathcal{T}} \int_{\mathcal{S}} |\theta(t, s)|^2 dt ds < \infty$ . Natural estimators of  $\mathbb{E}[\mathcal{X}]$  and  $\mathbb{E}[\mathcal{Y}]$  are the empirical means,  $\overline{\mathcal{X}} \triangleq 1/n \sum_{i=1}^n \mathcal{X}_i$  and  $\overline{\mathcal{Y}} \triangleq 1/n \sum_{i=1}^n \mathcal{Y}_i$ . The centered random error function  $\epsilon$  is assumed to be independent of  $\mathcal{X}$ , and  $\mathbb{E}\|\epsilon\|^2 < \infty$ .

We define the elements of the centered model as follows,  $X \triangleq \mathcal{X} - \mathbb{E}[\mathcal{X}]$  and  $Y \triangleq \mathcal{Y} - \mathbb{E}[\mathcal{Y}]$ , and the centered FFLRM writes

$$Y(s) = \int_{\mathcal{T}} X(t)\theta(t, s)dt + \epsilon(s). \quad (2.1)$$

Denote

$$C_X(t_1, t_2) = \mathbb{E}\left(X(t_1)X(t_2)\right), \quad t_1, t_2 \in \mathcal{T},$$

and

$$C_Y(s_1, s_2) = \mathbb{E}\left(Y(s_1)Y(s_2)\right), \quad s_1, s_2 \in \mathcal{S},$$

as the covariance functions of  $X$  and  $Y$ , respectively. As  $C_X \in \mathbb{L}^2(\mathcal{T}^2)$  and  $C_Y \in \mathbb{L}^2(\mathcal{S}^2)$  are a positive self-adjoint operators, then, according to Mercer's theorem ([Hsing and Eubank, 2015](#), Theorem 4.6.5),  $C_X$  and  $C_Y$  admit spectral expansions in  $\mathbb{L}^2(\mathcal{T}^2)$  and  $\mathbb{L}^2(\mathcal{S}^2)$  respectively

$$C_X(t_1, t_2) = \sum_{k=1}^{+\infty} \lambda_k \phi_k(t_1)\phi_k(t_2) \quad \text{and} \quad C_Y(s_1, s_2) = \sum_{j=1}^{+\infty} \mu_j \psi_j(s_1)\psi_j(s_2),$$

where  $\lambda_1 > \lambda_2 > \dots > 0$  and  $\mu_1 > \mu_2 > \dots > 0$  are the eigenvalue sequences of the covariance functions  $C_X$  and  $C_Y$ , respectively, while  $\{\phi_k\}_{k \geq 1}$  and  $\{\psi_j\}_{j \geq 1}$  are the corresponding orthonormal bases of eigenfunctions in  $\mathbb{L}^2(\mathcal{T})$  and  $\mathbb{L}^2(\mathcal{S})$ .

The Karhunen–Loève (KL) expansions of the curves  $X$  and  $Y$  (Hsing and Eubank, 2015, Theorem 7.3.5) in  $\mathbb{L}^2(\mathcal{T})$  and  $\mathbb{L}^2(\mathcal{S})$  are respectively

$$X(t) = \sum_{k=1}^{+\infty} \xi_k \phi_k(t) \quad \text{and} \quad Y(s) = \sum_{j=1}^{+\infty} \beta_j \psi_j(s),$$

for all  $t \in \mathcal{T}$  and  $s \in \mathcal{S}$ , where  $\xi_k = \int_{\mathcal{T}} X(t) \phi_k(t) dt$  and  $\beta_j = \int_{\mathcal{S}} Y(s) \psi_j(s) ds$  are uncorrelated random variables with zero mean and variances  $\mathbb{E}(\xi_k^2) = \lambda_k$  and  $\mathbb{E}(\beta_j^2) = \mu_j$  for all  $k, j \geq 1$ . These coefficients  $\xi_k$  and  $\beta_j$  are called functional principal components scores. By Parseval's identity and Fubini's theorem, we have

$$\sum_{k=1}^{+\infty} \mathbb{E}(\xi_k^2) < \infty \quad \text{and} \quad \sum_{j=1}^{+\infty} \mathbb{E}(\beta_j^2) < \infty.$$

Ramsay and Silverman (2005, Chapter 16, Section 1.1), Park and Qian (2012), Crambes and Mas (2013) and Imaizumi and Kato (2018) suggest that  $\theta(t, s)$  can be expressed in terms of a double basis expansion

$$\theta(t, s) = \sum_{k=1}^{+\infty} \sum_{j=1}^{+\infty} \theta_{k,j} \phi_k(t) \psi_j(s),$$

where  $\theta_{k,j} = \int_{\mathcal{T}} \int_{\mathcal{S}} \phi_k(t) \theta(t, s) \psi_j(s) dt ds$ .

Noticing that

$$\int_{\mathcal{T}} \theta(t, s) X(t) dt = \sum_{k=1}^{+\infty} \sum_{j=1}^{+\infty} \theta_{k,j} \xi_k \psi_j(s),$$

and

$$\mathbb{E}(\xi_k Y(s)) = \lambda_k \sum_{j=1}^{+\infty} \theta_{k,j} \psi_j(s),$$

we obtain the following characterization of the bivariate functional coefficient  $\theta(t, s)$  as

$$\theta(t, s) = \sum_{k=1}^{+\infty} \frac{\mathbb{E}(\xi_k Y(s))}{\lambda_k} \phi_k(t).$$

This characterization leads to a method for estimating  $\theta(t, s)$ . For example, Park and Qian (2012) and Crambes and Mas (2013) use the following truncation estimator with  $k_n \rightarrow \infty$  as  $n \rightarrow \infty$ ,

$$\hat{\theta}(t, s) = \frac{1}{n} \sum_{i=1}^n \sum_{k=1}^{k_n} \frac{\hat{\xi}_{ik} Y_i(s)}{\hat{\lambda}_k} \hat{\phi}_k(t), \quad (2.2)$$

where  $\hat{\lambda}_k$  and  $\hat{\phi}_k$  are the estimators of the eigenvalues  $\lambda_k$  and the eigenfunctions  $\phi_k$  and the estimates of the FPC scores are  $\hat{\xi}_{ik} = \int_{\mathcal{T}} X_i(t) \hat{\phi}_k(t) dt$ . We will use this characterization in this paper.

Benatia et al. (2017) propose an estimator similar to (2.2). The sum over  $k$  is not truncated, but regularized with the term  $\hat{\lambda}_k + \kappa$

$$\hat{\theta}^{\ddagger}(t, s) = \frac{1}{n} \sum_{i=1}^n \sum_{k=1}^n \frac{\hat{\xi}_{ik} Y_i(s)}{\hat{\lambda}_k + \kappa} \hat{\phi}_k(t).$$

Ramsay and Silverman (2005) obtain the following characterization of the bivariate functional coefficient, for some small or large numbers  $\iota_1$  and  $\iota_2$ , as

$$\theta^{\dagger}(t, s) = \sum_{k=1}^{\iota_1} \sum_{j=1}^{\iota_2} \theta_{k,j} \phi_k(t) \psi_j(s), \quad \text{where } \theta_{k,j} = \frac{\mathbb{E}(\xi_k \beta_j)}{\lambda_k}.$$

Therefore Imaizumi and Kato (2018) obtain the following characterization

$$\theta^{\dagger\dagger}(t, s) = \sum_{k=1}^{+\infty} \frac{\mathbb{E}(\xi_k \beta_j)}{\lambda_k} \phi_k(t) \psi_j(s).$$

Notice that  $\theta^{\dagger}(t, s)$  and  $\theta^{\dagger\dagger}(t, s)$  based on truncating the double series, namely, the double truncation, which will not be discussed here.

## 2.2 Operatorial point of view

We notice in this subsection that the model (2.1) can be seen from an operatorial point of view. Indeed, we can write the model

$$Y(\mathbf{s}) = (\Theta \cdot X)(\mathbf{s}) + \epsilon(\mathbf{s}),$$

where the slope operator  $\Theta : \mathbb{L}^2(\mathcal{T} \times \mathcal{S}) \rightarrow \mathbb{L}^2(\mathcal{S})$  is an integral continuous operator defined for all  $u \in \mathbb{L}^2(\mathcal{T})$  by  $(\Theta \cdot u)(s) = \int_{\mathcal{T}} u(t) \theta(t, s) dt$  which kernel is  $\theta \in \mathbb{L}^2(\mathcal{T} \times \mathcal{S})$ .

To close this section, we introduce the operators that will be used through the following theorems in which we describe prediction convergence rates. The covariance operator of  $X$ , denoted  $\Gamma$ , is defined by

$$\Gamma(u) = \mathbb{E}(X \otimes X(u)), \quad \text{for all } u \in \mathbb{L}^2(\mathcal{T}).$$

Note that the covariance operator is a natural extension of the covariance matrix, in the infinite dimensional framework. We also introduce the cross-covariance operator  $\Delta$  of  $(X, Y)$  given by

$$\Delta(u) = \mathbb{E}(Y \otimes X(u)), \quad \text{for all } u \in \mathbb{L}^2(\mathcal{T}).$$

For any integer  $k$ , we define  $\Pi_k$  the orthogonal projection operator on the subspace  $\text{Span}(\phi_1, \dots, \phi_k)$ , given by

$$\Pi_k = \sum_{j=1}^k \phi_j \otimes \phi_j.$$

Empirical counterparts of  $\Gamma$ ,  $\Delta$  and  $\Pi_k$ , respectively, are denoted  $\hat{\Gamma}_n$ ,  $\hat{\Delta}_n$  and  $\hat{\Pi}_{k_n}$ . These operators are naturally defined by  $\hat{\Gamma}_n = \frac{1}{n} \sum_{i=1}^n X_i \otimes X_i$ ,  $\hat{\Delta}_n = \frac{1}{n} \sum_{i=1}^n Y_i \otimes X_i$  and  $\hat{\Pi}_{k_n} = \sum_{j=1}^{k_n} \hat{\phi}_j \otimes \hat{\phi}_j$ . The functional principal component regression estimator  $\hat{\Theta}$  of  $\Theta$  is defined by

$$\hat{\Theta} = \hat{\Pi}_{k_n} \hat{\Delta}_n (\hat{\Pi}_{k_n} \hat{\Gamma}_n \hat{\Pi}_{k_n})^{-1}.$$

### 3 The centered Function-on-function linear model with partially observed covariate and response: Reconstructing $X$ and $Y$

In this section, we are interested in the most general case of missing data in the centered function-on-function linear regression: when both the functional covariate and the functional response are partially observed. We follow the methodology studied in [Kneip and Liebl \(2020\)](#) for reconstructing the missing parts of the curves. Next, once the initial sample is completed, we will present the estimation of the slope operator  $\Theta$  or its kernel  $\theta$  and predict new values for the response.

In the following, we denote "O<sup>[Y]</sup>", "M<sup>[Y]</sup>", "O<sup>[X]</sup>" and "M<sup>[X]</sup>" a given production of  $O_i^{[Y]}$ ,  $M_i^{[Y]}$ ,  $O_i^{[X]}$  and  $M_i^{[X]}$ , respectively corresponding to the observed domain and the missing domain of  $Y_i$  included in  $\mathcal{S}$  and the observed domain and the missing domain of  $X_i$  included in  $\mathcal{T}$ . In addition, the corresponding parts of the curve  $X_i$  are denoted  $X_i^{O^{[X]}}$  and  $X_i^{M^{[X]}}$ . Similarly, the corresponding parts of  $Y_i$  are denoted  $Y_i^{O^{[Y]}}$  and  $Y_i^{M^{[Y]}}$ .

#### 3.1 Curve reconstruction of the covariate and the response

The Karhunen–Loève expansions of the observed sample curves  $X_i^{O^{[X]}}$  in  $\mathbb{L}^2(O^{[X]})$  and  $Y_i^{O^{[Y]}}$  in  $\mathbb{L}^2(O^{[Y]})$  are written

$$X_i^{O^{[X]}}(t) = \sum_{k=1}^{+\infty} \xi_{ik}^{O^{[X]}} \phi_k^{O^{[X]}}(t) \quad \text{and} \quad Y_i^{O^{[Y]}}(s) = \sum_{j=1}^{+\infty} \beta_{ij}^{O^{[Y]}} \psi_j^{O^{[Y]}}(s),$$

where  $\xi_{ik}^{O^{[X]}} = \int_{\mathcal{T}} X_i^{O^{[X]}}(t) \phi_k^{O^{[X]}}(t) dt$  and  $\beta_{ij}^{O^{[Y]}} = \int_{\mathcal{S}} Y_i^{O^{[Y]}}(s) \psi_j^{O^{[Y]}}(s) ds$  are uncorrelated random variables with zero mean and variances  $\mathbb{E} \left( \xi_k^{O^{[X]}} \right)^2 = \lambda_k^{O^{[X]}}$  and  $\mathbb{E} \left( \beta_j^{O^{[Y]}} \right)^2 = \mu_j^{O^{[Y]}}$  for all  $k, j \geq 1$ .

We consider a reconstruction problem relating the missing part of the curves to the observed part, writing

$$X_i^{M^{[X]}}(t_2) = L^{[X]}(X_i^{O^{[X]}}(t_1)) + \mathcal{Z}_i^{[X]}(t_2), \quad \text{for all } t_1 \in O^{[X]} \text{ and } t_2 \in M^{[X]},$$

and

$$Y_i^{M^{[Y]}}(s_2) = L^{[Y]}(Y_i^{O^{[Y]}}(s_1)) + \mathcal{Z}_i^{[Y]}(s_2), \quad \text{for all } s_1 \in O^{[Y]} \text{ and } s_2 \in M^{[Y]},$$

where  $L^{[X]} : \mathbb{L}_2(O^{[X]}) \rightarrow \mathbb{L}_2(M^{[X]})$  and  $L^{[Y]} : \mathbb{L}_2(O^{[Y]}) \rightarrow \mathbb{L}_2(M^{[Y]})$  are linear reconstruction operators and  $\mathcal{X}_i^{[X]} \in \mathbb{L}_2(M^{[X]})$  and  $\mathcal{X}_i^{[Y]} \in \mathbb{L}_2(M^{[Y]})$  are reconstruction errors. We need to minimize the mean square error between the curves with fragmentary data and the linear reconstruction operators  $L^{[X]}$  and  $L^{[Y]}$  as follows

$$\mathbb{E}\left(\|X_i^{M^{[X]}} - L^{[X]}(X_i^{O^{[X]}})\|^2\right) \quad \text{and} \quad \mathbb{E}\left(\|Y_i^{M^{[Y]}} - L^{[Y]}(Y_i^{O^{[Y]}})\|^2\right). \quad (3.1)$$

Thes operator  $L^{[X]}$  (or  $L^{[Y]}$ ) has been studied by [Kraus \(2015\)](#) and [Kneip and Liebl \(2020\)](#). [Kraus \(2015\)](#) proposed to use the ridge regularization method. The estimator is introduced by  $\mathcal{L}^{(\alpha)} = \mathcal{R}_{MO}\mathcal{R}_{OO}^{(\alpha)-1}$  where  $\mathcal{R}$  is the covariance operator defined on  $\mathbb{L}^2(\mathcal{T})$  and  $\mathcal{R}_{OO}^{(\alpha)-1} = \mathcal{R}_{OO} + \alpha \mathcal{I}_O$  where  $\alpha$  is a positive parameter and  $\mathcal{I}_O$  is the identity operator on  $\mathbb{L}^2(O^{[X]})$ . Besides, [Kraus and Stefanucci \(2020\)](#) prove that the reconstruction operator (considered in [Kraus \(2015\)](#)) can be seen as Hilbert-Schmidt integral operator from  $\mathbb{L}^2(O^{[X]})$  to  $\mathbb{L}^2(O^{[X]})$  writing

$$L(X_i^{O^{[X]}}) = \int_{O^{[X]}} a_i(\cdot, t) X_i^{O^{[X]}}(t) dt,$$

for all  $i = 1, \dots, n$ , where  $a$  is a square integrable function on  $M^{[X]} \times O^{[X]}$ .

The optimal reconstruction operators minimizing (3.1) are denoted  $\mathcal{L}(X_i^{O^{[X]}})$  and  $\mathcal{J}(Y_i^{O^{[Y]}})$  and defined by

$$\mathcal{L}(X_i^{O^{[X]}})(t_2) = \sum_{k=1}^{+\infty} \xi_{ik}^{O^{[X]}} \frac{\langle \phi_k^{O^{[X]}}, \gamma_{t_2} \rangle}{\lambda_k^{O^{[X]}}} \quad \text{and} \quad \mathcal{J}(Y_i^{O^{[Y]}})(s_2) = \sum_{j=1}^{+\infty} \beta_{ij}^{O^{[Y]}} \frac{\langle \psi_j^{O^{[Y]}}, \gamma_{s_2} \rangle}{\mu_j^{O^{[Y]}}}, \quad (3.2)$$

where  $\gamma_{t_2}(t_1) = \mathbb{E}\left(X_i^{M^{[X]}}(t_2) X_i^{O^{[X]}}(t_1)\right)$ , for all  $t_1 \in O^{[X]}$  and  $t_2 \in M^{[X]}$ , and  $\gamma_{s_2}(s_1) = \mathbb{E}\left(Y_i^{M^{[Y]}}(s_2) Y_i^{O^{[Y]}}(s_1)\right)$ , for all  $s_1 \in O^{[Y]}$  and  $s_2 \in M^{[Y]}$ .

The operators in relation (3.2) are estimated as in [Kneip and Liebl \(2020, Section 2\)](#) and [Crambes et al. \(2023, Subsection 2.1 and Subsection 2.2\)](#) by  $\widehat{\mathcal{L}}_{k_n}(X_i^{O^{[X]}})$  and  $\widehat{\mathcal{J}}_{j_n}(Y_i^{O^{[Y]}})$ , where the truncation parameters  $k_n$  and  $j_n$  are positive integers that can be fixed automatically with a grid search. The solution of (3.2) uses local linear smoothers for unknown quantities in (3.2), considering the following notations. Let  $\kappa_1$  and  $\kappa'_1$  be kernels and  $h_X$  and  $h_Y$  be bandwidths of the local linear smoothers of the curves  $X_i^{O^{[X]}}$  and  $Y_i^{O^{[Y]}}$ , respectively. Moreover, let  $\kappa_2$  and  $\kappa'_2$  be bivariate kernels and  $h_{\gamma_{t_2}}$  and  $h_{\gamma_{s_2}}$  be bandwidths of the local linear smoothers of the covariance functions  $\gamma_{t_2}$  and  $\gamma_{s_2}$ , respectively.

In the following, we consider the whole sample  $\mathcal{D}_n \triangleq \{(X_1^*, Y_1^*), \dots, (X_n^*, Y_n^*)\}$ , with possibly reconstructed curves on the missing parts, that is

$$X_i^*(t) = \begin{cases} X_i^{O^{[X]}}(t) & \text{if } t \in O^{[X]}, \\ \widehat{\mathcal{L}}_{k_n}(X_i^{O^{[X]}})(t) & \text{if } t \in M^{[X]} \end{cases} \quad \text{and} \quad Y_i^*(s) = \begin{cases} Y_i^{O^{[Y]}}(s) & \text{if } s \in O^{[Y]}, \\ \widehat{\mathcal{J}}_{j_n}(Y_i^{O^{[Y]}})(s) & \text{if } s \in M^{[Y]}. \end{cases}$$



### 3.2 Estimation of slope operator and its kernel and prediction

We estimate the kernel function  $\theta$  with

$$\hat{\theta}^*(t, s) = \frac{1}{n} \sum_{i=1}^n \sum_{k=1}^{k_n} \frac{\hat{\xi}_{ik,rec}^* Y_i^*(s)}{\hat{\lambda}_{k,rec}^*} \hat{\phi}_{k,rec}^*(t), \quad (3.3)$$

where  $\hat{\phi}_{1,rec}^*, \dots, \hat{\phi}_{k_n,rec}^*$  and  $\hat{\lambda}_{1,rec}^*, \dots, \hat{\lambda}_{k_n,rec}^*$  represent respectively the  $k_n$  first eigenfunctions and eigenvalues of the covariance operator  $\hat{\Gamma}_{n,rec}^*$  and  $\hat{\xi}_{ik,rec}^* = \int_{\mathcal{S}} X_i^*(t) \hat{\phi}_{k,rec}^*(t) dt$  are the estimates of the FPC scores.

From an operatorial point of view, the covariance operator  $\Gamma_{rec}^*$  of  $X^*$  and the cross-covariance operator  $\Delta_{rec}^*$  of  $(X^*, Y^*)$  are given by  $\Gamma_{rec}^* = \mathbb{E}[X^* \otimes X^*]$  and  $\Delta_{rec}^* = \mathbb{E}[Y^* \otimes X^*]$ , for all  $u \in \mathbb{L}^2(\mathcal{S})$ . The orthogonal projection operator  $\Pi_{k,rec}^*$  on the subspace  $\text{Span}(\phi_{1,rec}^*, \dots, \phi_{k,rec}^*)$  is given by  $\Pi_k^* = \sum_{j=1}^k \phi_{j,rec}^* \otimes \phi_{j,rec}^*$ . The empirical counterparts of  $\Gamma_{rec}^*$ ,  $\Delta_{rec}^*$  and  $\Pi_{k,rec}^*$ , are denoted respectively  $\hat{\Gamma}_{n,rec}^*$ ,  $\hat{\Delta}_{n,rec}^*$  and  $\hat{\Pi}_{k_n,rec}^*$ . The estimator of the slope operator  $\Theta$  is given by

$$\hat{\Theta}^* = \langle \hat{\theta}^*(\cdot, s), \cdot \rangle = \hat{\Pi}_{k_n,rec}^* \hat{\Delta}_{n,rec}^* \left( \hat{\Pi}_{k_n,rec}^* \hat{\Gamma}_{n,rec}^* \hat{\Pi}_{k_n,rec}^* \right)^{-1}.$$

Finally, we obtain the prediction of the response when a new explanatory curve  $X_{new}$  is given, by

$$Y_{new}^*(s) = \int_{\mathcal{S}} X_{new}^*(t) \hat{\theta}^*(t, s) dt = \frac{1}{n} \sum_{i=1}^n \sum_{k=1}^{k_n} \frac{\hat{\xi}_{ik,rec}^* \hat{\xi}_{new;k,rec}^* Y_i^*(s)}{\hat{\lambda}_{k,rec}^*},$$

for all  $s \in \mathcal{S}$ , where  $\hat{\xi}_{new;k,rec}^* = \int_{\mathcal{S}} X_{new}^*(t) \hat{\phi}_{k,rec}^*(t) dt$ . Alternatively, it is also given with the operatorial quantities

$$Y_{new}^* = \hat{\Pi}_{k_n,rec}^* \hat{\Delta}_{n,rec}^* \left( \hat{\Pi}_{k_n,rec}^* \hat{\Gamma}_{n,rec}^* \hat{\Pi}_{k_n,rec}^* \right)^{-1} X_{new}^*.$$

### 3.3 Assumptions

To achieve our theoretical results, we need first to adopt some classical assumptions which have been similarly used in [Kneip and Liebl \(2020\)](#) and [Crambes et al. \(2023\)](#) to control the curve reconstruction for the covariate and the response.

**(A.1)**  $X$  and  $Y$  have finite fourth moment order.

**(A.2)** Let  $np \rightarrow \infty$  when  $n \rightarrow \infty$ , where  $p = p(n)$  is the number of observation points of the covariate. Similarly,  $nq \rightarrow \infty$  when  $n \rightarrow \infty$ , where  $q = q(n)$  is the number of observation points of the response. We assume in the following that  $p = n^{\eta_1}$  with  $0 < \eta_1 < \infty$  and  $q = n^{\zeta_1}$  with  $0 < \zeta_1 < \infty$ .

**(A.3)** • The bandwidth  $h_X$  satisfies  $h_X \rightarrow 0$  and  $(ph_X) \rightarrow \infty$  as  $p \rightarrow \infty$ . For instance, we assume that  $h_X = \frac{1}{n^{\eta_2}}$  with  $0 < \eta_2 < \eta_1$ . The bandwidth  $h_{\gamma_{t_2}}$  satisfies  $h_{\gamma_{t_2}} \rightarrow 0$  and  $(n(p^2 - p)h_{\gamma_{t_2}}) \rightarrow \infty$  as  $n(p^2 - p) \rightarrow \infty$ . For example, we can take  $h_{\gamma_{t_2}} = \frac{1}{n^{\eta_3}}$  with

$0 < \eta_3 < 2\eta_1 + 1$ .

- Let  $\kappa_1$  and  $\kappa_2$  be nonnegative, second order univariate and bivariate kernel functions with support  $[-1, 1]$ . For example, we can use univariate and bivariate Epanechnikov kernel functions with compact support  $[-1, 1]$ , namely  $\kappa_1(x) = \frac{3}{4}(1-x^2)\mathbb{1}_{[-1,1]}(x)$  and  $\kappa_2(x, y) = \frac{9}{16}(1-x^2)(1-y^2)\mathbb{1}_{[-1,1]}(x)\mathbb{1}_{[-1,1]}(y)$ .

**(A.4)** • For any subinterval  $O^{[X]} \subseteq \mathcal{T}$ , we assume that the eigenvalues  $\lambda_1 > \lambda_2 > \dots > 0$  have multiplicity one. Moreover, we assume that there exist  $a_O > 1$  and  $0 < c_O < \infty$  such that (i)  $\lambda_k^{O^{[X]}} - \lambda_{k+1}^{O^{[X]}} \geq c_O k^{-a_O-1}$ , (ii)  $\lambda_k^{O^{[X]}} = \mathcal{O}(k^{-a_O})$ , (iii)  $1/\lambda_k^{O^{[X]}} = \mathcal{O}(k^{a_O})$  as  $k \rightarrow \infty$ .

- $\mathbb{E}(\xi_k^4) = \mathcal{O}(\lambda_k^2)$ .
- For any subinterval  $O^{[X]} \subseteq \mathcal{T}$ , we assume that there exists  $0 < A_O < \infty$  such that the eigenfunctions satisfy  $\sup_{t \in \mathcal{T}} \sup_{k \geq 1} |\tilde{\phi}_k^{O^{[X]}}(t)| \leq A_O$ , where  $\tilde{\phi}_k^{O^{[X]}}(t_2) = \langle \phi_k^{O^{[X]}}, \gamma_{t_2} \rangle / \lambda_k^{O^{[X]}}$ .

**(A.5)** • The bandwidth  $h_Y$  satisfies  $h_Y \rightarrow 0$  and  $(qh_Y) \rightarrow \infty$  as  $q \rightarrow \infty$ . Moreover, we assume that  $h_Y = \frac{1}{n^{\zeta_2}}$  with  $0 < \zeta_2 < \zeta_1$ . The bandwidth  $h_{\gamma_{s_2}}$  satisfies  $h_{\gamma_{s_2}} \rightarrow 0$  and  $(n(q^2 - q)h_{\gamma_{s_2}}) \rightarrow \infty$  as  $n(q^2 - q) \rightarrow \infty$ . For example, we can take  $h_{\gamma_{s_2}} = \frac{1}{n^{\zeta_3}}$  with  $0 < \zeta_3 < 2\zeta_1 + 1$ .

- Let  $\kappa'_1$  and  $\kappa'_2$  be nonnegative, second order univariate and bivariate kernel functions with support  $[-1, 1]$ .

**(A.6)** • For any subinterval  $O^{[Y]} \subseteq \mathcal{S}$ , we assume that  $\mu_1 > \mu_2 > \dots > 0$  have multiplicity one and we assume that there exist  $b_O > 1$  and  $0 < d_O < \infty$  such that (j)  $\mu_j^{O^{[Y]}} - \mu_{j+1}^{O^{[Y]}} \geq d_O j^{-b_O-1}$ , (jj)  $\mu_j^{O^{[Y]}} = \mathcal{O}(j^{-b_O})$ , (jjj)  $1/\mu_j^{O^{[Y]}} = \mathcal{O}(j^{b_O})$  as  $j \rightarrow \infty$ .

- $\mathbb{E}(\beta_k^4) = \mathcal{O}(\mu_k^2)$ .
- For any subinterval  $O^{[Y]} \subseteq \mathcal{S}$ , we assume that there exists  $0 < B_O < \infty$  such that  $\sup_{s \in \mathcal{S}} \sup_{j \geq 1} |\tilde{\psi}_j^{O^{[Y]}}(s)| \leq B_O$ , where  $\tilde{\psi}_j^{O^{[Y]}}(s_2) = \langle \psi_j^{O^{[Y]}}, \gamma_{s_2} \rangle / \lambda_j^{O^{[Y]}}$ .

Assumption **(A.1)** holds for many processes  $X$  and  $Y$  (Gaussian processes, bounded processes). Assumption **(A.2)** is mild and can be satisfied even if the number of observation points  $p$  and  $q$  do not go fast to infinity. **(A.3)** and **(A.5)** are classic assumptions in the context of local polynomials smoothers. Assumptions **(A.4)** and **(A.6)** are similar to the ones in [Kneip and Liebl \(2020\)](#).

### 3.4 Asymptotic results

Under assumptions **(A.1)**-**(A.6)**, it is proved in [Kneip and Liebl \(2020\)](#) that, in the case where  $p \sim n^{\eta_1}$  and  $q \sim n^{\zeta_1}$  with  $\eta_1, \zeta_1 \leq 1/2$  we have for any  $t \in \mathcal{T}$  and  $s \in \mathcal{S}$

$$|X_i^*(t) - X_i(t)| = \mathcal{O}_p\left(n^{-\eta_1(a_O-1)/(2(a_O+2))}\right) \text{ and } |Y_i^*(s) - Y_i(s)| = \mathcal{O}_p\left(n^{-\zeta_1(b_O-1)/(2(b_O+2))}\right).$$

The previous result allows to obtain some bounds between quantities related to functional principal components analysis between the reconstructed curves and with the original curves.

We finish this subsection with the main result giving a bound for the prediction error of  $Y_{new}$  with a new value of the covariate  $X_{new}$ .

**Theorem 3.1.** *Under assumptions (A.1)-(A.6), taking  $k_n \sim p^{1/(a_O+2)}$ ,  $p \sim n^{\eta_1}$ ,  $j_n \sim q^{1/(b_O+2)}$  and  $q \sim n^{\zeta_1}$ , with  $\eta_1, \zeta_1 \leq 1/2$ , we get*

$$\mathbb{E} \left( \left\| \left( \hat{\Theta}^* \cdot X_{new}^* - \Theta \cdot X_{new}^* \right) \right\|^2 \right) = \mathcal{O} \left( n^{-\eta_1(a_O-1)/(2(a_O+2))} + n^{\eta_1/(a_O+2)-1-\zeta_1(b_O-1)/(b_O+2)} \right).$$

**Corollary 3.2.** *We make a comparison between the parameters to find the best convergence error. We summarize the error rates in Table 1.*

Table 1: Convergence error rates depending on the observation points and the regularity of the curves  $X$  and  $Y$ .

(i) $\eta_1 = \zeta_1$	$a_O \leq b_O$	$\mathcal{O} \left( n^{-\eta_1(a_O-1)/(2(a_O+2))} \right)$
	$a_O > b_O$	$\mathcal{O} \left( n^{\eta_1/(a_O+2)-1-\eta_1(b_O-1)/(b_O+2)} \right)$
(ii) $\eta_1 < \zeta_1$		$\mathcal{O} \left( n^{\eta_1/(a_O+2)-1-\zeta_1(b_O-1)/(b_O+2)} \right)$
(iii) $\eta_1 > \zeta_1$		$\mathcal{O} \left( n^{-\eta_1(a_O-1)/(2(a_O+2))} \right)$

Comparing the parameters, we remark all the convergence rates depend in particular on the parameter  $a_O > 1$ , which is directly linked to the smoothness of the stochastic process  $X$ . The larger  $a_O$  is, the smoother  $X$  is. In the case (i) when the number of observation points of the covariate is equivalent to that of the response for  $a_O > b_O$  and in the case (ii), the convergence rates depend of the parameter  $a_O$  and also of the parameter  $b_O > 1$ , which is directly linked to the smoothness of the stochastic process  $Y$ . In these cases, the final rate of convergence will be linked to the parameter  $a_O$  or  $b_O$  corresponding to the less smooth process (either  $X$  or  $Y$ ).

## 4 The centered function-on-function with partially observed covariate and response: Reconstructing $X$ and imputing $Y$

We have seen in the previous section the methodology for reconstructing the missing parts of the explanatory curves. In this section, we try another strategy to deal with missing data on the response. After reconstructing the missing parts of the covariate  $X$ , we apply the regression imputation methodology as presented in Crambes et al. (2023) for a real response. Next, we will present the estimation of the kernel function  $\theta$  and predict the new response once all sample is completed.

#### 4.1 Regression imputation on the functional response

We consider the following missing data mechanism for the response, through a variable  $\delta^{[Y]}$  leading to the sample  $(\delta_i^{[Y]})_{i=1,\dots,n}$  such that,

$$\delta_i^{[Y]} = \begin{cases} 0 & \text{if } M_i^{[Y]} \neq \emptyset, \\ 1 & \text{if } O_i^{[Y]} = \mathcal{S}. \end{cases}$$

We assume that the response is missing at random (MAR), which means that the fact that  $Y$  contains missing parts does not depend on the response of the model, but can possibly depend on the reconstructed covariate,

$$\mathbb{P}(\delta^{[Y]} = 1 \mid X^*, Y) = \mathbb{P}(\delta^{[Y]} = 1 \mid X^*).$$

We denote the number of curves partially observed by

$$m_n^{[Y]} = \sum_{i=1}^n \mathbf{1}_{\{\delta_i^{[Y]}=0\}}.$$

Using the exponent notation "obs" to make reference to the units for which the response is completely observed, we define the covariance operator  $\Gamma_{rec}^{obs}$  and the cross-covariance operator  $\Delta_{rec}^{obs}$  with the reconstructed curves by  $\Gamma_{rec}^{obs} = \mathbb{E}(\delta^{[Y]} X^* \otimes X^*)$  and  $\Delta_{rec}^{obs} = \mathbb{E}(\delta^{[Y]} Y \otimes X^*)$ . The empirical counterparts of  $\Gamma_{rec}^{obs}$  and  $\Delta_{rec}^{obs}$  are denoted respectively  $\hat{\Gamma}_{n,rec}^{obs}$  and  $\hat{\Delta}_{n,rec}^{obs}$ .

Let  $Y_\ell$  be a response curve such that  $\delta_\ell^{[Y]} = 0$ , we define the imputed response  $Y_{\ell,imp}$  by

$$Y_{\ell,imp}(s) = \int_{\mathcal{T}} X_\ell^*(t) \tilde{\theta}(t, s) dt,$$

with

$$\tilde{\theta}(t, s) = \frac{1}{n - m_n^{[Y]}} \sum_{i=1}^n \sum_{k=1}^{k_n} \frac{\hat{\xi}_{ik,rec}^{obs} \delta_i^{[Y]}(s) Y_i^*(s)}{\hat{\lambda}_{k,rec}^{obs}} \hat{\phi}_{k,rec}^{obs}(t), \quad (4.1)$$

for all  $(t, s) \in \mathcal{T} \times \mathcal{S}$ , where  $\hat{\xi}_{ik,rec}^{obs} = \int_{\mathcal{T}} X_i^*(t) \hat{\phi}_{k,rec}^{obs}(t) dt$  and  $\hat{\phi}_{1,rec}^{obs}, \dots, \hat{\phi}_{k_n,rec}^{obs}$  and  $\hat{\lambda}_{1,rec}^{obs}, \dots, \hat{\lambda}_{k_n,rec}^{obs}$  represent respectively the  $k_n$  first eigenfunctions and eigenvalues of the covariance operator  $\hat{\Gamma}_{n,rec}^{obs}$ .

Alternatively, if we denote  $\hat{\Pi}_{k_n,rec}^{obs}$  the projection on the space spanned by the  $k_n$  first eigenfunctions of  $\hat{\Gamma}_{n,rec}^{obs}$ , the estimation of the slope operator  $\Theta$  is given by

$$\tilde{\Theta} = \hat{\Pi}_{k_n,rec}^{obs} \hat{\Delta}_{n,rec}^{obs} \left( \hat{\Pi}_{k_n,rec}^{obs} \hat{\Gamma}_{n,rec}^{obs} \hat{\Pi}_{k_n,rec}^{obs} \right)^{-1}.$$

and the imputation  $Y_{\ell,imp}$  can also be written

$$Y_{\ell,imp}(s) = \hat{\Pi}_{k_n,rec}^{obs} \hat{\Delta}_{n,rec}^{obs} \left( \hat{\Pi}_{k_n,rec}^{obs} \hat{\Gamma}_{n,rec}^{obs} \hat{\Pi}_{k_n,rec}^{obs} \right)^{-1} X_\ell^*(s),$$

## 4.2 Estimation of the slope operator and its kernel and prediction

Once the whole database has been reconstructed, we estimate the bivariate functional coefficient  $\theta$  with

$$\hat{\theta}^{**}(t, s) = \frac{1}{n} \sum_{i=1}^n \sum_{k=1}^{k_n} \frac{\hat{\xi}_{ik,rec}^* Y_i^{**}(s)}{\hat{\lambda}_{k,rec}^*} \hat{\phi}_{k,rec}^*(t), \quad (4.2)$$

for all  $(t, s) \in \mathcal{T} \times \mathcal{S}$ , where  $Y_i^{**} = Y_i \delta_i^{[Y]} + Y_{i,imp}(1 - \delta_i^{[Y]})$  for all  $i = 1, \dots, n$ . The estimation of the operator  $\Theta$  is similarly given by

$$\hat{\Theta}^{**} = \hat{\Pi}_{k_n,rec}^* \hat{\Delta}_{n,rec}^{**} \left( \hat{\Pi}_{k_n,rec}^* \hat{\Gamma}_{n,rec}^* \hat{\Pi}_{k_n,rec}^* \right)^{-1},$$

where  $\hat{\Delta}_{n,rec}^{**}$  is the empirical counterpart of the cross-covariance operator  $\Delta_{rec}^{**}$ .

We use this estimation to predict a new value of the response  $Y$  when a new explanatory curve  $X_{new}$  is given by

$$\hat{Y}_{new}(s) = \int_{\mathcal{T}} X_{new}^*(t) \hat{\theta}^{**}(t, s) d(t),$$

which can also be written

$$\hat{Y}_{new} = \hat{\Pi}_{k_n,rec}^* \hat{\Delta}_{n,rec}^{**} \left( \hat{\Pi}_{k_n,rec}^* \hat{\Gamma}_{n,rec}^* \hat{\Pi}_{k_n,rec}^* \right)^{-1} X_{new}^*.$$

## 4.3 Asymptotic results

The proof of these results follows the same lines as the proof of **Theorem (3.1)** and **Theorem (3.3)** in [Crambes et al. \(2023\)](#).

**Theorem 4.1.** *Under assumptions (A.1)-(A.4), if we take  $k_n \sim p^{1/(a_O+2)}$  and  $p \sim n^{\eta_1}$  with  $\eta_1 \leq 1/2$ , we obtain*

$$\mathbb{E} \left( \left\| \left( \tilde{\Theta} \cdot X_{\ell}^* - \Theta \cdot X_{\ell}^* \right) \right\|^2 \right) = \mathcal{O} \left( n^{-\eta_1(a_O-1)/(2(a_O+2))} + \frac{n^{\eta_1/(a_O+2)}}{n - m_n^{[Y]}} \right),$$

for  $\ell \in \tilde{\mathcal{D}}_m$ , where  $\tilde{\mathcal{D}}_m$  is the set of missing responses of size  $m_n^{[Y]}$ .

**Theorem 4.2.** *Under assumptions (A.1)-(A.4), and  $k_n \sim p^{1/(a_O+2)}$  and  $p \sim n^{\eta_1}$  with  $\eta_1 \leq 1/2$ , the prediction error is*

$$\mathbb{E} \left( \left\| \left( \hat{\Theta}^{**} \cdot X_{new}^* - \Theta \cdot X_{new}^* \right) \right\|^2 \right) = \mathcal{O} \left( n^{-\eta_1(a_O-1)/(2(a_O+2))} + \frac{n^{\eta_1/(a_O+2)}}{n - m_n^{[Y]}} \right).$$

**Remark 4.3.** *Comparing parameters as  $n^{\eta_1/(a_O+2)-1-\zeta_1(b_O-1)/(b_O+2)} \lesssim \frac{n^{\eta_1/(a_O+2)}}{n - m_n^{[Y]}}$ , we find that the prediction error with reconstruction (obtained in Section 3) is asymptotically at least the same than the prediction error with imputation.*

**Remark 4.4.** Theoretical results are generally obtained under assumptions concerning the rate of convergence of the integer  $k_n$ . In practice, this integer is selected by minimizing a certain empirical criterion such as Generalized Cross Validation (GCV) criterion, Cross Validation (CV) criterion, or  $K$ -fold Cross Validation ( $K$ -fold CV) criterion. We chose in the following simulation section the GCV procedure, known to be computationally fast. The GCV criteria is given as follows for imputation

$$GCV(k_n) = \frac{(n - m_n^{[Y]}) \sum_{i=1}^n \left\| \left( \tilde{\Theta} \cdot X_i^* - \Theta \cdot X_i^* \right) \right\|^2 \delta_i^{[Y]}}{\left( (n - m_n^{[Y]}) - k_n \right)^2},$$

and analogously for prediction.

## 5 Simulations

### 5.1 Methodology

In this section, we conduct Monte Carlo experiments to illustrate the finite-sample performance of the proposed methods presented in section 3 and section 4. We set  $\mathcal{S} = \mathcal{T} = [0, 1]$ . Each response and predictor curve is observed at  $q = 90$  and  $p = 100$  equally spaced points in their domains, respectively. For computational simplicity we consider equidistant points  $s_j = j/(q - 1)$ ,  $j = 0, \dots, q - 1$ , and  $t_k = k/(p - 1)$ ,  $k = 0, \dots, p - 1$ . Two simulated data mechanisms are generated. Each model is defined by

$$Y^{(w)}(s) = \int_0^1 X^{(w)}(t) \theta^{(w)}(t, s) dt + \epsilon^{(w)}(s), \quad (5.1)$$

for  $t \in \mathcal{T}$  and  $s \in \mathcal{S}$ , where  $w = 1, 2$ . We approximate the integral in (5.1) using a Riemann sum over the grid  $\mathbf{t}$ . The analytical expressions of the kernels and processes are given below.

SCENARIO 1 The kernel is given by

$$\theta^{(1)}(t, s) = \sum_{k=1}^{50} \sum_{j=1}^{50} \theta_{k,j} \phi_k(t) \psi_j(s),$$

where  $\phi_k(t) = \sqrt{2} \cos(k\pi t)$ ,  $\psi_j(s) = \sqrt{2} \cos(j\pi s)$  and  $\theta_{k,j} = 4(-1)^{k+j} k^{-2.5} j^{-3}$ . The input is the random function  $X^{(1)}(t) = \sum_{k=1}^{50} k^{-1} \xi_k \phi_k(t)$ , where the  $\xi_k$ 's are independently sampled from the uniform distribution on  $[-\sqrt{3}, \sqrt{3}]$ . Finally, the noise is given by  $\epsilon^{(1)} = \sum_{j=1}^{50} \beta'_j \psi_j(s)$  and the  $\beta'_j$ 's are independent mean-zero Gaussian random variables with variances equal to  $j^{-1.1}$ . Then, the covariance function  $C_{X^{(1)}}$  of  $X^{(1)}$  is given by

$$C_{X^{(1)}}(t_1, t_2) = \sum_{k=1}^{50} \frac{2}{k^2} \cos(k\pi t_1) \cos(k\pi t_2), \quad \text{where } t_1, t_2 \in \mathcal{T}.$$

SCENARIO 2 The kernel is defined by  $\theta^{(2)}(t, s) = s^3 + \sin(2\pi t)^3$  and the noise  $\epsilon^{(2)}$  is generated according to a standard Brownian motion divided by 20. In addition, the functional covariate  $X^{(2)}$  is generated through its covariance function, defined, for all  $t_1, t_2 \in \mathcal{T}$ , by

$$C_{X^{(2)}}(t_1, t_2) = \frac{\sigma_1^2 \exp(-|t_1 - t_2|^\alpha)}{\zeta},$$

with  $\sigma_1 = 1$ ,  $\alpha = 2$  and  $\zeta = 0.2$ . In this setting, even if a polynomial decrease of the eigenvalues of the covariance operator of  $X^{(2)}$  is required in our theoretical results (see assumption **(A.4)**), we want to see how the method works in practice if this assumption is no more satisfied, namely here in the case of an exponential decay.

Figure 1 shows 10 discretized predictor functions  $X_i^{(1)}$ , the error functions  $\epsilon_i^{(1)}$  and the response functions  $Y_i^{(1)}$ . Figure 2 shows 10 discretized predictor functions  $X_i^{(2)}$ , the error functions  $\epsilon_i^{(2)}$  and the response functions  $Y_i^{(2)}$ .

Figure 3 shows the covariance functions of the covariate for SCENARIO 1 and 2. Figure 4 shows kernel functions for SCENARIO 1 and 2.

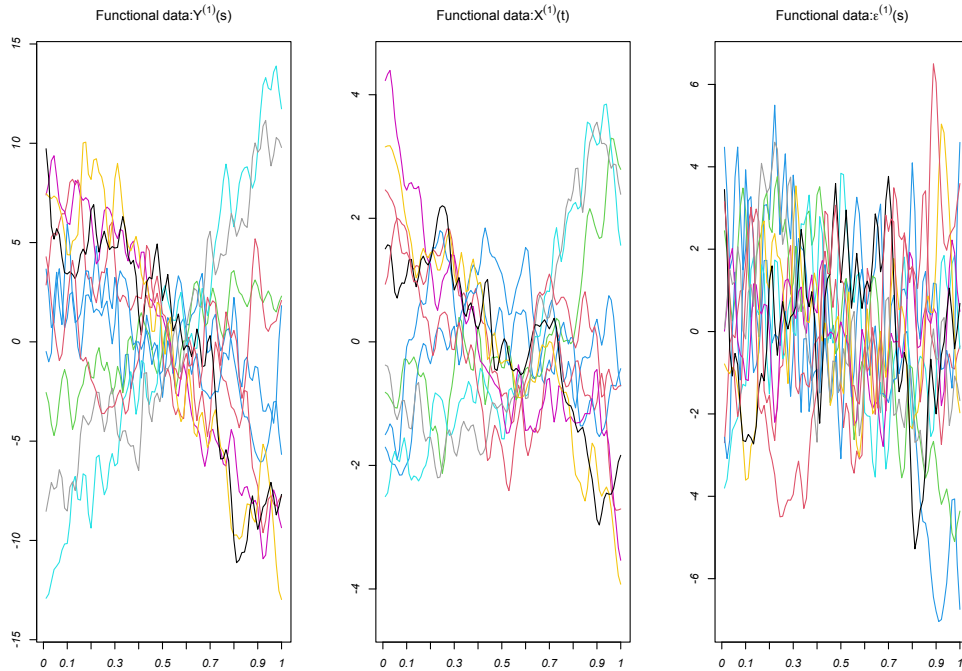


Figure 1: Examples of simulated functions with SCENARIO 1.

To deal with partially observed curves for the covariate and response, we adopted the missing data simulation scenario from [Kneip and Liebl \(2020\)](#) and [Crambes et al. \(2023\)](#) such that

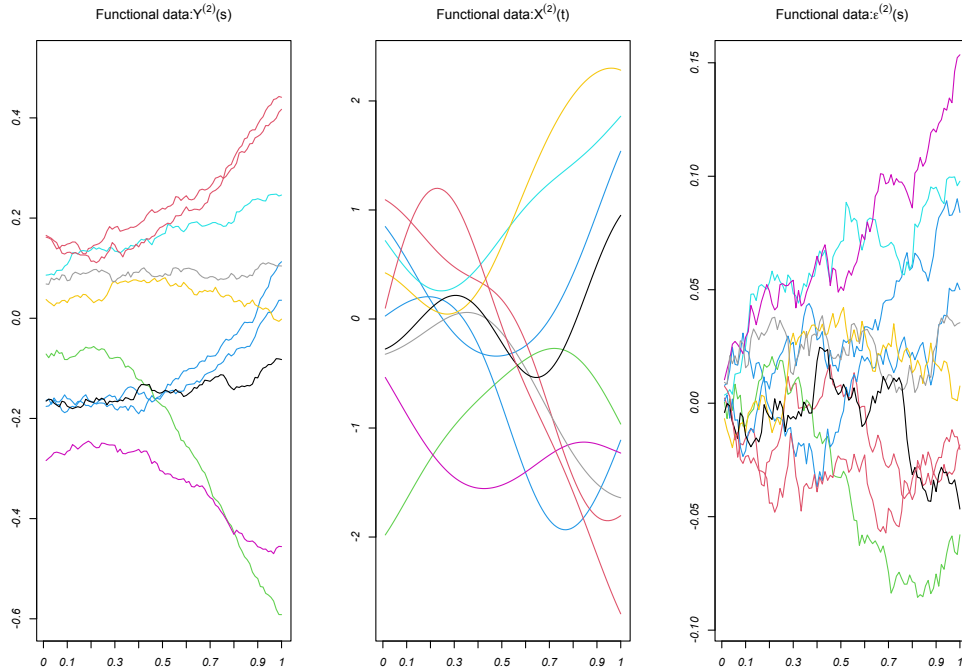


Figure 2: Examples of simulated functions with SCENARIO 2.

- 70% (respectively 55%) of the curves are fully observed on  $[0, 1]$ ,
- for the 30% (respectively 45%) of partially observed curves, the curves  $X_i$  and  $Y_i$  are fully observed on  $[A_i, B_i] \subset [0, 1]$  with  $A_i$  drawn with uniform law on the interval  $[0, A]$  and  $B_i = A_i + B$ , with either  $A = 1/50$  and  $B = 49/50$  or  $A = 3/50$  and  $B = 47/50$  for SCENARIO 1. We take either  $A = 1/50$  and  $B = 49/50$  or  $A = 5/50$  and  $B = 45/50$  for SCENARIO 2.

We simulate the number of missing data on the response  $Y$  and the indicator  $\delta^{[Y]}$  by the logistic functional regression. The variable  $\delta$  follows the Bernoulli law with parameter  $p(X)$  such that

$$\log\left(\frac{p(X)}{1-p(X)}\right) = \left[ \int_S \left( \int_{\mathcal{T}} |s-t| X(t) dt \right)^2 ds \right]^{1/2} + ct,$$

where  $ct$  is a constant allowing to take different levels of missing data. For exemple,  $ct = 2$  gives around 9.727% of missing data and  $ct = 0.2$  gives around 38.801% of missing data.

Let us notice that we use a spline smoothed version of the different estimators (2.2), (3.3), (4.1) and (4.2), according to the so-called Smooth Principal Components Regression (SPCR) from Cardot et al. (2003). Let us remark that, with appropriate conditions on the spline parameters, all the theoretical results obtained in our work will also apply when using the SPCR



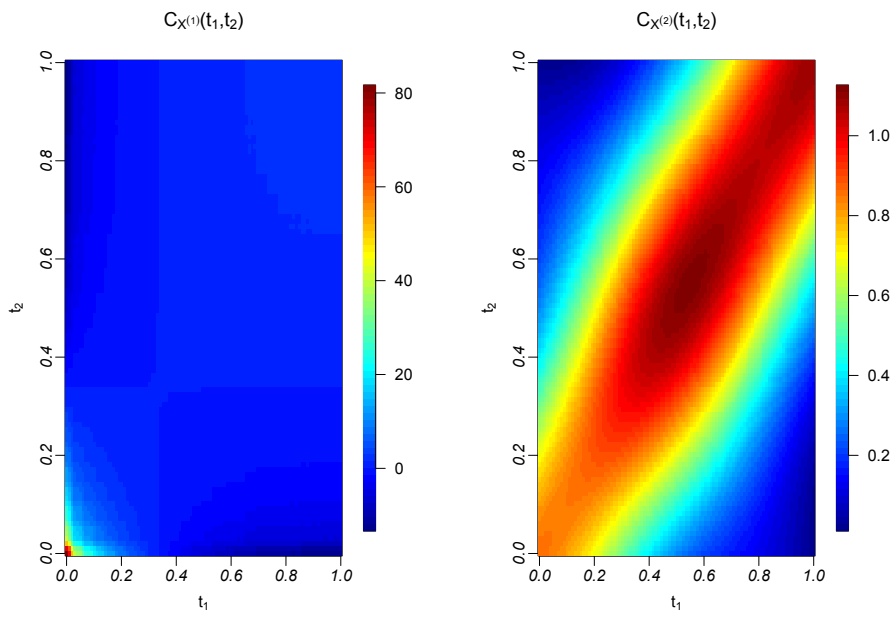


Figure 3: The covariance functions for SCENARIO 1 and 2.

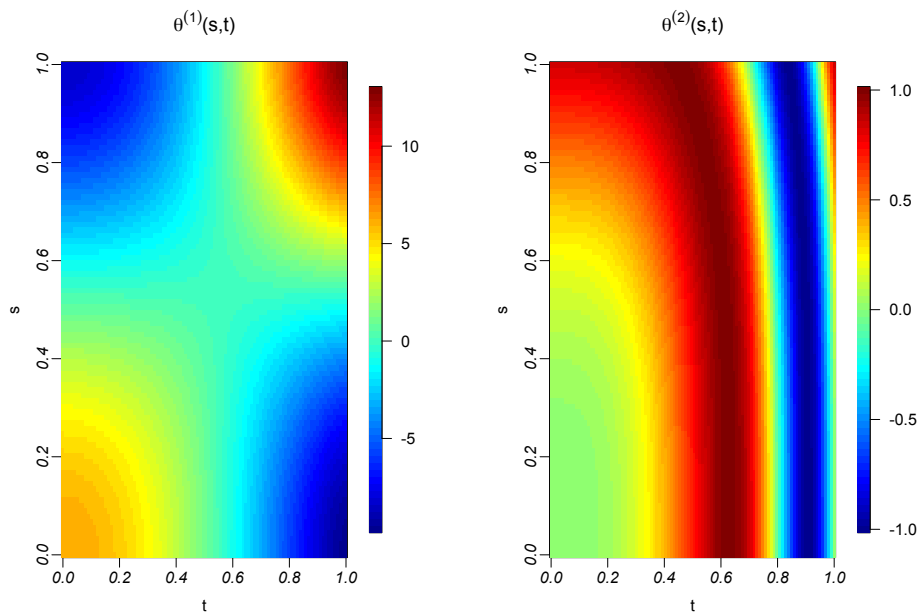


Figure 4: The kernel functions for SCENARIO 1 and 2.

estimation. We use a regression spline basis with 20 knots, a degree 3 and the order of derivation 2. The choice of these parameters is not crucial in our study, especially in comparison with the choice of the number of principal components. The choice of this optimal tuning parameter is made on a growing sequence of dimension  $k_n = 2, \dots, 22$ .

The dataset of size  $N$  is randomly splitted into a training set of size  $n = \frac{4}{5}N$  and a test set of size  $n_1 = \frac{1}{5}N$ . We consider sample sizes  $N = 500, 1300$ . For each scenario, we use 200 Monte Carlo runs for the model assessment. In all numerical experiments, the proposed estimators have been carried out with the free software R.

## 5.2 Criteria

We use the following criteria to evaluate the performance of the methods.

- Criterion 1:  $\overline{MSPE} = \frac{1}{Sim} \sum_{j=1}^{Sim} MSPE(j)$  is the average mean square prediction error. This criterion tends to zero when the sample size tends to infinity, where  $MSPE(j) = \frac{1}{n_1} \sum_{\ell=n+1}^{n+n_1} \left\| \left( \hat{\Theta} \cdot X_{\ell}^{*,j} - \Theta \cdot X_{\ell}^{*,j} \right) \right\|^2$  is the mean square prediction error computed on the  $j^{th}$  simulated sample,  $j \in \{1, \dots, Sim\}$ .
- Criterion 2:  $\overline{RT} = \frac{1}{Sim} \sum_{j=1}^{Sim} RT(j)$  is the average ratio respect to truth. This criterion tends to one when the sample size tends to infinity, where  $RT(j) = \frac{\sum_{\ell=n+1}^{n+n_1} \left\| \left( \hat{\Theta} \cdot X_{\ell}^{*,j} \right) - Y_{\ell}^j \right\|^2}{\sum_{\ell=n+1}^{n+n_1} \left\| e_{\ell}^j \right\|^2}$  is the ratio between the mean square prediction error and the mean square prediction error when the true parameters are known, computed on the  $j^{th}$  simulated sample.

We consider another criterion which is the determination coefficient  $R^2$ . In this context of functional regression setting, several definitions exist. Given the fitted values  $\hat{Y}_i(s)$ , we used the definition as in [Harezlak et al. \(2007\)](#) given by

$$R^2 = \frac{1}{|\mathcal{S}|} \int_{\mathcal{S}} R^2(s) ds = \frac{1}{|\mathcal{S}|} \int_{\mathcal{S}} \left( 1 - \frac{\sum_{i=1}^n \left( Y_i(s) - \hat{Y}_i(s) \right)^2}{\sum_{i=1}^n Y_i(s)^2} \right) ds.$$

## 5.3 Simulation results

We denote the methods presented in this paper by :

- **Reconst\_X\_Y** :  $X$  and  $Y$  are partially observed , the missing parts of  $X$  and  $Y$  are reconstructed.
- **Reconst\_X, Imp\_Y** :  $X$  and  $Y$  are partially observed, the missing parts of  $X$  are reconstructed and  $Y$  imputed.

Moreover, we compare to other methods :

- **Full\_X\_Y** :  $X$  and  $Y$  are fully observed, this corresponds to the complete reference dataset.
- **Reconst\_X, Remov\_Y** :  $X$  and  $Y$  are partially observed, the missing parts of  $X$  are reconstructed and the missing part of  $Y$  are removed from the sample.
- **Reconst\_X, Remov\_Y** :  $X$  and  $Y$  are partially observed, the individuals presenting partially observed curves are removed from the sample.

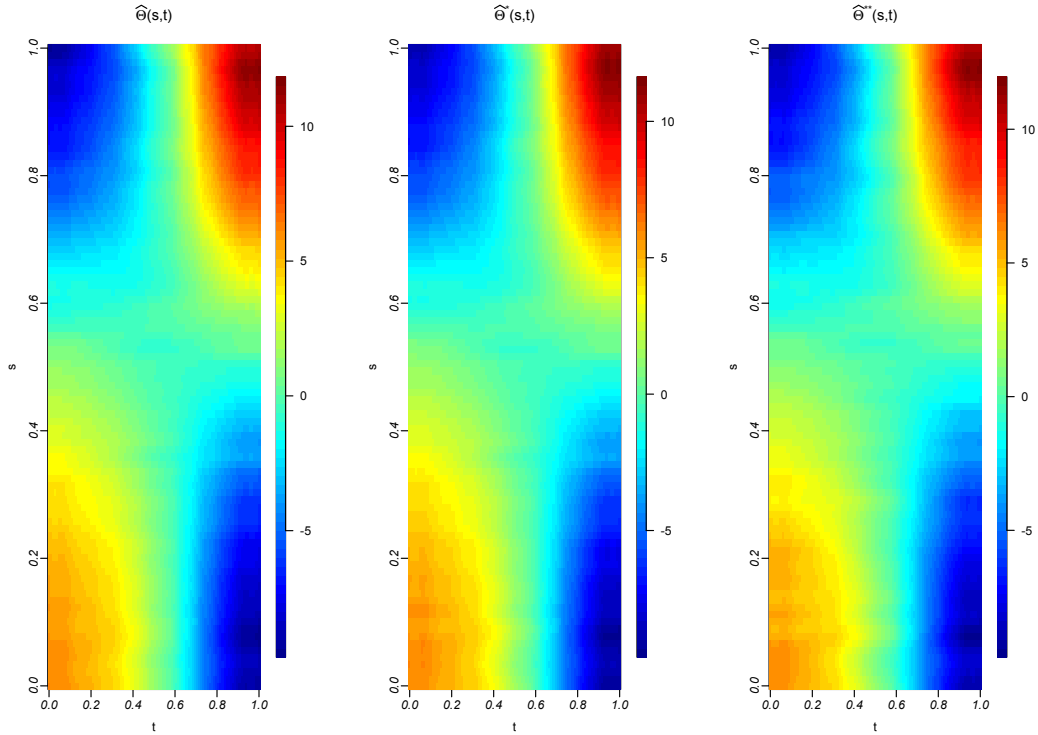


Figure 5: The estimated coefficient functions for SCENARIO 1.

Even if our main goal is prediction, Figure 5 show estimates of the kernel function in SCENARIO 1 (with a sample size  $N = 400$ ) for the dimension  $k_n^*$  chosen by the GCV criterion, respectively with full data ( $\hat{\Theta}$ ), reconstruction of the missing parts of  $X$  and  $Y$  ( $\hat{\Theta}^*$ ) and reconstruction of the missing parts of  $X$  and imputation of  $Y$  ( $\hat{\Theta}^{**}$ ). The missing part is 12% for both curves  $X$  and  $Y$ , the observed part being  $[3/50, 47/50]$ . Moreover, 39.375% of curves  $Y$  (with  $ct = 0.1$ ) are affected by missing data and 42.250% of curves  $X$  are affected by missing data. We remark that the estimators look graphically quite close,  $\hat{\Theta}^*$  seems to be a little closer to  $\hat{\Theta}$  than  $\hat{\Theta}^{**}$ . A similar plot is obtained for SCENARIO 2 (Figure 6) with 37.812% of curves  $Y$  (with  $ct = 0.1$ ) affected by missing data and 46.750% of curves  $X$  affected by missing data. In this situation,  $\hat{\Theta}^*$  seems much closer to  $\hat{\Theta}$  than  $\hat{\Theta}^{**}$ .

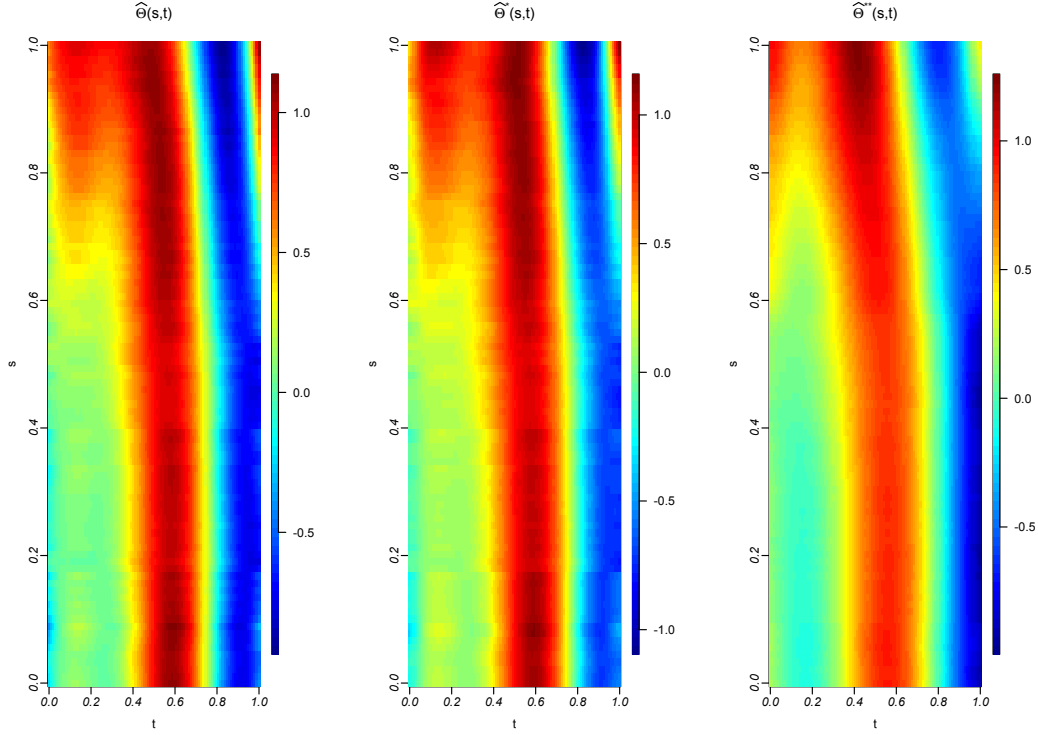


Figure 6: The estimated coefficient functions for SCENARIO 2.

We give in Table 7 the values of the determination coefficient  $R^2$  and the value  $k_n^*$  chosen by the GCV criterion both scenarios 1 and 2. In scenario 1, we get a worse  $R^2$  coefficient, maybe due to the fact that the curves  $X$  are not so smooth and do not seem easy to reconstruct.

Table 2:  $R^2$  and  $k_n^*$  for scenarios 1 and 2.

Methods		SCENARIO 1	SCENARIO 2
<b>Full_X_Y</b>	$R^2$	68.996 %	98.652 %
	$k_n^*$	2	6
<b>Reconst_X_Y</b>	$R^2$	68.859 %	98.641 %
	$k_n^*$	2	6
<b>Reconst_X, Imp_Y</b>	$R^2$	68.853 %	98.630 %
	$k_n^*$	2	5

Tables 3, 4, 5 and 6 give the values of the criteria  $\overline{MSPE}$  and  $\overline{RT}$  for scenarios 1 and 2 with different values of sample size, and different levels of missing data. The first conclusion is the fact that the errors decrease as the sample size increases. Secondly, these errors increase with the percentage of missing data on  $X$  or on  $Y$ . The rate of missing data on  $Y$  seems to

have a more important impact on the errors, whatever the scenario we consider. In all cases, the method **Reconst\_X\_Y** reconstructing both curves  $X$  and  $Y$  has a better behaviour than the method **Reconst\_X, Imp\_Y** reconstructing  $X$  and imputing  $Y$ , which is quite in accordance to our theoretical results. The part of the observed curve is an important parameter in the curve reconstruction: as it can be expected, the results are better when the curve reconstruction is easier (for example when the observed part is  $[1/50, 49/50]$ , corresponding to 4% of missing information on the curves). Results tend to deteriorate when the curve reconstruction is harder (for example when the observed part is  $[3/50, 47/50]$ , corresponding to 12% of missing information on the curves). Finally, these two methods behave better than the other more naive methods (**Reconst\_X, Remov\_Y** and **Reconst\_X, Remov\_Y**) that partially or completely ignore missing individuals affected by missing data.

Table 3: Mean and standard deviation errors for the predicted values based on 200 simulation replications with different levels of missing data and a sample size 500 (left panel) and a sample size 1300 (right panel). Partially observed curves are fully observed on [1/50, 49/50] with SCENARIO 1.

Rate of missing data in Y in %	10.150 (1.407)	10.181 (1.481)	40.599 (2.742)	40.542 (2.596)
Rate of missing data in X in %	30.114 (1.998)	44.814 (2.498)	29.778 (2.354)	44.958 (2.180)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	27.618 (7.073)	27.472 (7.162)	27.620 (7.584)	27.312 (7.077)
$\overline{RT} \times 10$	10.073 (0.066)	10.067 (0.058)	10.071 (0.057)	10.081 (0.066)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	28.159 (7.149)	28.234 (7.270)	28.843 (8.222)	28.804 (7.654)
$\overline{RT} \times 10$	10.075 (0.068)	10.069 (0.059)	10.072 (0.059)	10.084 (0.070)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^3</math></b>	30.388 (7.754)	30.547 (8.388)	43.315 (13.043)	41.853 (10.940)
$\overline{RT} \times 10$	10.080 (0.070)	10.074 (0.061)	10.113 (0.073)	10.120 (0.082)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^3</math></b>	30.588 (8.032)	30.609 (8.464)	43.981 (13.870)	42.239 (11.563)
$\overline{RT} \times 10$	10.080 (0.071)	10.074 (0.061)	10.114 (0.075)	10.120 (0.084)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	40.575 (11.757)	48.537 (14.961)	58.408 (19.486)	70.458 (21.764)
$\overline{RT} \times 10$	10.105 (0.101)	10.112 (0.108)	10.147 (0.113)	10.198 (0.138)

Rate of missing data in Y in %	10.174 (0.991)	10.236 (0.933)	40.345 (1.694)	40.293 (1.659)
Rate of missing data in X in %	30.085 (1.238)	44.919 (1.352)	30.030 (1.326)	44.911 (1.391)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	12.418 (2.692)	12.473 (3.016)	12.686 (3.498)	12.164 (2.636)
$\overline{RT} \times 10$	10.029 (0.026)	10.031 (0.027)	10.030 (0.026)	10.033 (0.025)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	12.909 (2.726)	13.231 (3.013)	13.906 (3.982)	13.399 (3.101)
$\overline{RT} \times 10$	10.030 (0.027)	10.034 (0.028)	10.033 (0.027)	10.036 (0.028)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^3</math></b>	14.110 (3.114)	14.111 (3.178)	19.008 (4.963)	19.135 (4.669)
$\overline{RT} \times 10$	10.032 (0.026)	10.036 (0.029)	10.047 (0.031)	10.052 (0.035)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^3</math></b>	14.138 (3.103)	14.152 (3.227)	19.538 (5.113)	19.401 (4.769)
$\overline{RT} \times 10$	10.032 (0.026)	10.036 (0.029)	10.049 (0.032)	10.052 (0.035)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	18.444 (4.846)	22.823 (5.951)	25.462 (6.564)	30.748 (8.185)
$\overline{RT} \times 10$	10.047 (0.036)	10.058 (0.045)	10.061 (0.044)	10.077 (0.068)

Table 4: Mean and standard deviation errors for the predicted values based on 200 simulation replications with different levels of missing data and a sample size 500 (left panel) and a sample size 1300 (right panel). Partially observed curves are fully observed on [3/50, 47/50] with SCENARIO 1.

Rate of missing data in $Y$ in %	10.250 (1.502)	10.193 (1.607)	40.324 (2.412)	40.250 (2.415)
Rate of missing data in $X$ in %	29.915 (2.010)	44.870 (2.301)	29.830 (2.067)	44.879 (2.156)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	26.989 (7.726)	27.259 (7.507)	28.020 (7.759)	27.176 (7.059)
$\overline{RT} \times 10$	10.072 (0.063)	10.069 (0.060)	10.078 (0.063)	10.068 (0.062)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	34.408 (8.040)	38.372 (8.323)	36.244 (8.611)	38.411 (7.747)
$\overline{RT} \times 10$	10.089 (0.078)	10.098 (0.081)	10.098 (0.077)	10.095 (0.083)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^3</math></b>	36.752 (8.608)	40.348 (8.839)	49.018 (10.995)	51.103 (11.675)
$\overline{RT} \times 10$	10.094 (0.081)	10.105 (0.082)	10.123 (0.088)	10.127 (0.094)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^3</math></b>	36.891 (8.717)	40.403 (8.860)	49.436 (11.610)	51.639 (12.131)
$\overline{RT} \times 10$	10.095 (0.082)	10.105 (0.083)	10.124 (0.087)	10.127 (0.094)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	39.619 (10.117)	47.940 (14.026)	56.390 (16.332)	69.935 (21.684)
$\overline{RT} \times 10$	10.100 (0.087)	10.125 (0.111)	10.147 (0.111)	10.174 (0.118)
Rate of missing data in $Y$ in %	10.174 (0.965)	10.033 (0.898)	40.423 (1.426)	40.514 (1.548)
Rate of missing data in $X$ in %	30.182 (1.232)	45.107 (1.459)	30.108 (1.285)	44.896 (1.290)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	12.573 (3.071)	12.141 (2.684)	12.608 (3.125)	12.303 (2.797)
$\overline{RT} \times 10$	10.034 (0.026)	10.031 (0.027)	10.033 (0.027)	10.034 (0.029)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	19.898 (3.654)	22.796 (3.328)	20.396 (3.815)	23.408 (3.604)
$\overline{RT} \times 10$	10.052 (0.038)	10.060 (0.044)	10.053 (0.041)	10.057 (0.045)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^3</math></b>	21.153 (3.999)	23.828 (3.659)	26.195 (5.075)	28.732 (5.266)
$\overline{RT} \times 10$	10.053 (0.040)	10.063 (0.044)	10.067 (0.045)	10.068 (0.049)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^3</math></b>	21.180 (3.969)	23.847 (3.694)	26.477 (5.229)	29.101 (5.365)
$\overline{RT} \times 10$	10.054 (0.040)	10.063 (0.044)	10.068 (0.046)	10.069 (0.049)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^3</math></b>	18.892 (4.792)	22.105 (5.681)	25.893 (6.873)	31.192 (8.474)
$\overline{RT} \times 10$	10.047 (0.042)	10.058 (0.051)	10.067 (0.050)	10.073 (0.063)

Table 5: Mean and standard deviation errors for the predicted values based on 200 simulation replications with different levels of missing data and a sample size 500 (left panel) and a sample size 1300 (right panel). Partially observed curves are fully observed on [1/50, 49/50] with SCENARIO 2.

Rate of missing data in $Y$ in %	9.727	9.757	38.801	38.865	9.726	9.678	39.232	39.113
	(1.575)	(1.456)	(2.297)	(2.383)	(0.947)	(0.974)	(1.551)	(1.417)
Rate of missing data in $X$ in %	29.826	44.816	30.032	45.152	29.847	45.090	30.013	45.147
	(2.069)	(2.192)	(2.253)	(2.288)	(1.290)	(1.365)	(1.203)	(1.425)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	20.654	19.593	20.815	19.787	7.646	7.919	8.067	7.931
	(10.200)	(8.722)	(9.723)	(9.110)	(3.042)	(3.148)	(3.528)	(3.736)
$\overline{RT} \times 10$	10.158	10.143	10.194	10.153	10.062	10.065	10.065	10.055
	(0.219)	(0.210)	(0.234)	(0.202)	(0.085)	(0.086)	(0.079)	(0.082)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	20.706	19.614	22.114	20.838	7.740	8.009	8.917	8.860
	(10.140)	(8.743)	(10.572)	(9.366)	(3.070)	(3.224)	(3.640)	(3.916)
$\overline{RT} \times 10$	10.158	10.142	10.208	10.162	10.062	10.066	10.072	10.063
	(0.219)	(0.211)	(0.241)	(0.209)	(0.085)	(0.088)	(0.082)	(0.087)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^6</math></b>	22.473	21.336	32.038	31.248	8.441	8.809	12.872	12.691
	(11.216)	(9.245)	(16.620)	(15.914)	(3.599)	(3.633)	(5.396)	(6.048)
$\overline{RT} \times 10$	10.170	10.159	10.291	10.240	10.067	10.072	10.103	10.098
	(0.229)	(0.228)	(0.298)	(0.254)	(0.091)	(0.096)	(0.101)	(0.121)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^6</math></b>	22.473	21.340	32.052	31.312	8.442	8.809	12.938	12.702
	(11.213)	(9.239)	(16.651)	(15.948)	(3.599)	(3.632)	(5.390)	(6.052)
$\overline{RT} \times 10$	10.171	10.159	10.291	10.241	10.067	10.072	10.104	10.098
	(0.229)	(0.228)	(0.298)	(0.256)	(0.099)	(0.101)	(0.101)	(0.012)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	30.380	38.660	46.638	58.802	12.175	15.342	17.406	23.005
	(13.412)	(19.485)	(21.421)	(30.753)	(5.109)	(6.675)	(7.836)	(10.587)
$\overline{RT} \times 10$	10.206	10.293	10.403	10.518	10.106	10.107	10.146	10.175
	(0.322)	(0.438)	(0.451)	(0.577)	(0.138)	(0.154)	(0.144)	(0.189)



Table 6: Mean and standard deviation errors for the predicted values based on 200 simulation replications with different levels of missing data and a sample size 500 (left panel) and a sample size 1300 (right panel). Partially observed curves are fully observed on [5/50, 45/50] with SCENARIO 2.

Rate of missing data in $Y$ in %	9.520 (1.629)	9.684 (1.552)	39.322 (2.519)	38.958 (2.277)
Rate of missing data in $X$ in %	30.069 (2.096)	45.063 (2.232)	30.075 (2.106)	45.245 (2.213)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	19.781 (9.412)	20.747 (9.748)	20.282 (10.321)	19.100 (8.074)
$\overline{RT} \times 10$	10.169 (0.217)	10.170 (0.229)	10.176 (0.222)	10.160 (0.217)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	24.236 (10.292)	26.528 (10.650)	32.528 (12.805)	32.342 (11.271)
$\overline{RT} \times 10$	10.198 (0.222)	10.204 (0.240)	10.274 (0.261)	10.271 (0.258)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^6</math></b>	24.832 (10.676)	27.026 (11.140)	34.701 (14.806)	35.151 (14.391)
$\overline{RT} \times 10$	10.209 (0.235)	10.203 (0.249)	10.288 (0.295)	10.281 (0.281)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^6</math></b>	24.845 (10.685)	27.032 (11.154)	34.788 (14.898)	35.305 (14.574)
$\overline{RT} \times 10$	10.209 (0.235)	10.205 (0.249)	10.289 (0.295)	10.281 (0.282)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	29.948 (13.803)	39.318 (18.900)	44.115 (20.894)	56.488 (31.158)
$\overline{RT} \times 10$	10.259 (0.346)	10.345 (0.387)	10.390 (0.430)	10.458 (0.548)

Rate of missing data in $Y$ in %	9.629 (0.963)	9.673 (0.946)	39.176 (1.550)	39.195 (1.502)
Rate of missing data in $X$ in %	29.976 (1.284)	44.939 (1.488)	29.942 (1.325)	44.967 (1.392)
<b>Full_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	7.923 (3.677)	7.563 (3.144)	7.556 (3.433)	7.916 (3.391)
$\overline{RT} \times 10$	10.056 (0.082)	10.065 (0.079)	10.063 (0.087)	10.067 (0.078)
<b>Reconst_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	10.834 (3.977)	11.297 (3.315)	16.881 (5.605)	17.037 (4.891)
$\overline{RT} \times 10$	10.082 (0.088)	10.096 (0.086)	10.132 (0.117)	10.139 (0.109)
<b>Reconst_X, Imp_Y : <math>\overline{MSPE} \times 10^6</math></b>	11.062 (4.226)	11.409 (3.421)	14.049 (5.293)	15.743 (5.268)
$\overline{RT} \times 10$	10.082 (0.094)	10.097 (0.091)	10.118 (0.107)	10.127 (0.117)
<b>Reconst_X, Remov_Y : <math>\overline{MSPE} \times 10^6</math></b>	11.067 (4.222)	11.412 (3.423)	14.086 (5.289)	15.773 (5.290)
$\overline{RT} \times 10$	10.082 (0.094)	10.098 (0.091)	10.119 (0.108)	10.128 (0.118)
<b>Remov_X_Y : <math>\overline{MSPE} \times 10^6</math></b>	12.702 (5.567)	15.481 (7.721)	16.638 (7.140)	22.266 (11.556)
$\overline{RT} \times 10$	10.092 (0.120)	10.125 (0.156)	10.151 (0.150)	10.195 (0.224)

## 6 Real dataset study: Hawaii Ocean data

The Hawaii ocean time-series program has been making repeated observations of various hydrographic, chemical and biological properties of the water column at a station north of Oahu, Hawaii since October, 1988. In the CTD dataset<sup>1</sup> of this program, various variables were measured every two meters between 0 and 200 meters below the sea surface. These variables are viewed as functions of depth. The layer between 0 and 200 meters below the sea surface is called the epipelagic zone (or sunlight zone), where enough light is available for photosynthesis. Therefore, in this zone, the primary production in the ocean occurs, and plants and animals are largely concentrated. The measurements were repeated at different dates, which we consider as different sample curves. We use three functional variables: *Temperature*, *Salinity* and *Oxygen*. Each curve being observed at 101 equally spaced points in  $[0, 200]$ , we use the Oxygen as the response curve, expressed as a function of temperature, and we use the Salinity as the explanatory variable, also expressed as a function of temperature. We study the relationship between Salinity and Oxygen with the following model, written in the operatorial point of view:

$$Oxygen_i(Temperature) = \left( \Theta \cdot Salinity_i \right) (Temperature) + \varepsilon_i(Temperature), \quad i = 1, \dots, n, \quad \text{with } n = 191.$$

The graphical display of the initial sample of 191 pairs curves  $\{Oxygen_i, Salinity_i\}_{i=1}^{191}$  (raw curves and smooth curves) can be observed in Figure 7. We can see that, on these data, all the curves are partially observed, which is a different situation compared to the simulations realized in the previous section, where only a percentage of curves were partially observed. In particular, the imputation method presented in Section 4 cannot be directly applied. We want to explore in this situation a possibility of a hybrid use of both reconstruction and imputation of the response curves.

We consider two scenarios for the response curves and for each scenario we use two samples: a training sample of size  $\ell_1 = 153$  equal to  $4/5$  of the initial sample from which the estimates are computed and a testing sample of size  $\ell_2 = 38$  equal to  $1/5$  of the initial sample on which the prediction errors are calculated. First, we reconstruct the missing part of all covariates using the observed parts. Then, as all the response curves are partially observed, we cannot directly use the regression imputation method from Section 4. As a consequence, we choose to reconstruct only a certain percentage of the response curves.

- 'Scenario 1': we reconstruct more response curves (87.958% of initial response curves which is equivalent to 168 response curves) using the observed parts while leaving curves with a low percentage of missing part.
- 'Scenario 2': we reconstruct less response curves (47.644% of initial response curves which is equivalent to 91 response curves) using the observed parts while leaving curves with a high percentage of missing part.

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<sup>1</sup><https://hahana.soest.hawaii.edu/hot/hot-dogs/cextraction.html>

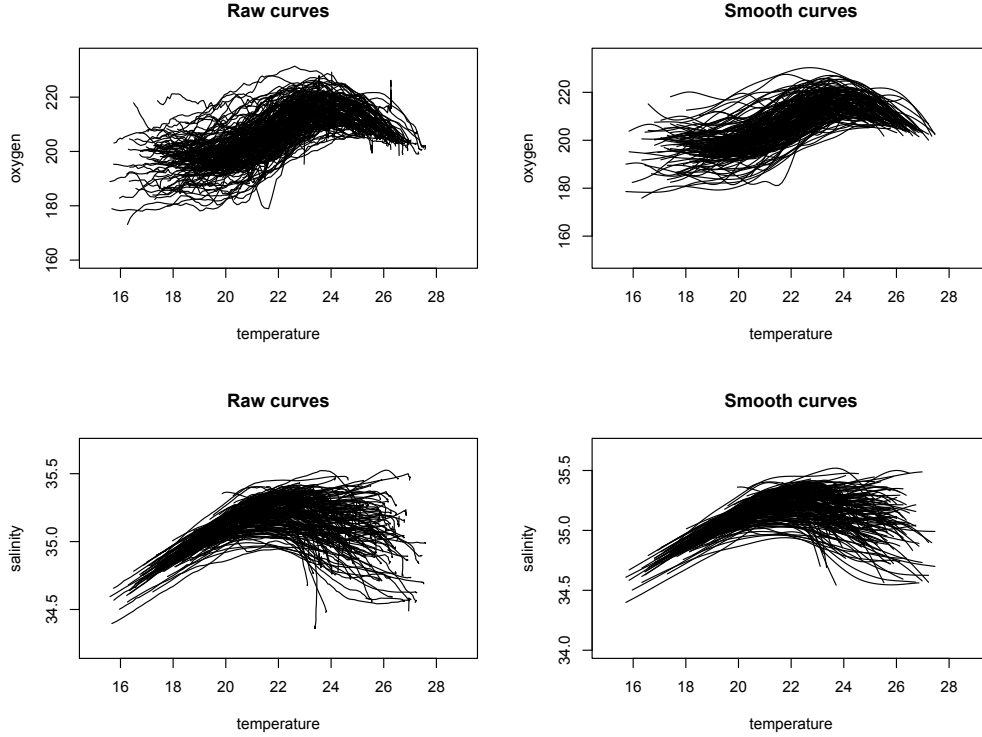


Figure 7: Partially Observed Functions: Raw curves of the oxygen variable (top left) and the salinity variable (bottom left). Smooth curves of the oxygen variable (top right) and the salinity variable (bottom right).

For 'Scenario 1', the training sample contains 23 partially observed response curves and 130 observed response curves and for 'Scenario 2' the training sample contains 100 partially observed response curves and 53 observed response curves. Moreover, for 'Scenario 1', 12.042% of response curves are affected by missing data. For 'Scenario 2', 52.356% of response curves are affected by missing data.

The graphical display of the training (resp. testing) samples of 153 (resp. 38) pairs of curves  $\{Oxygen_i, Salinity_i\}_{i=1}^{153(38)}$  can be observed in Figures 8 and 9.

We consider the MSPE criterion as in Section 5 to evaluate the quality of prediction error. We also consider a MSPA (Mean Square Prediction Absolute error) criterion, defined in the same way, replacing the 2-norm with the 1-norm in the definition of MSPE. Table 7 gives MSPE and MSPA criteria for both 'Scenario 1' and 'Scenario 2'. We can observe that the prediction errors are much better in 'Scenario 1'. In other words, in such situations where all the response curves are partially observed, it seems better to reconstruct a more important part of them, and then use the imputation method on the remaining part.

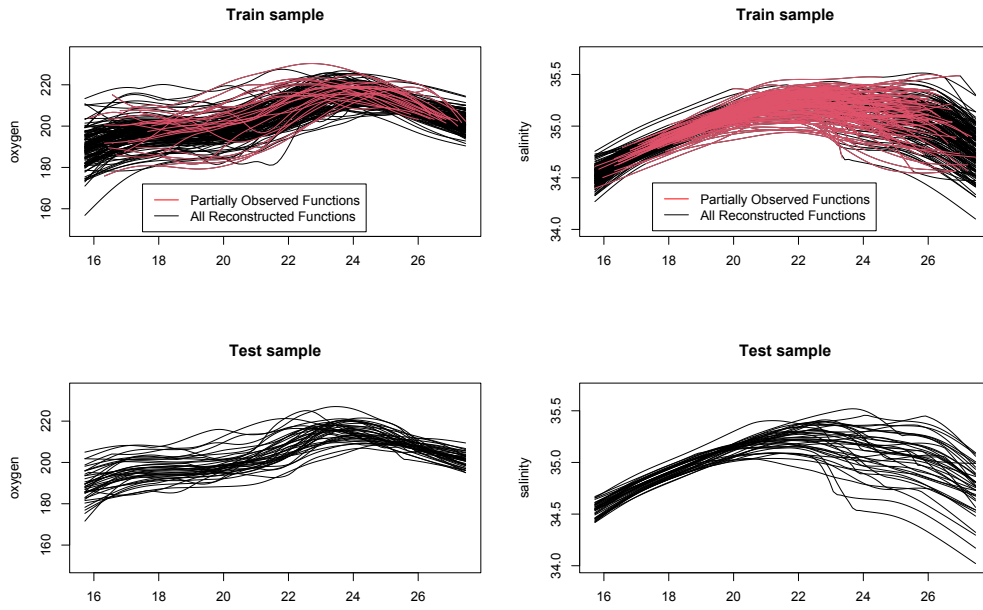


Figure 8: Training (resp. testing) samples of 153 (resp. 38) pairs of curves  $\{Oxygen_i, Salinity_i\}_{i=1}^{153(38)}$  in 'Scenario 1'.

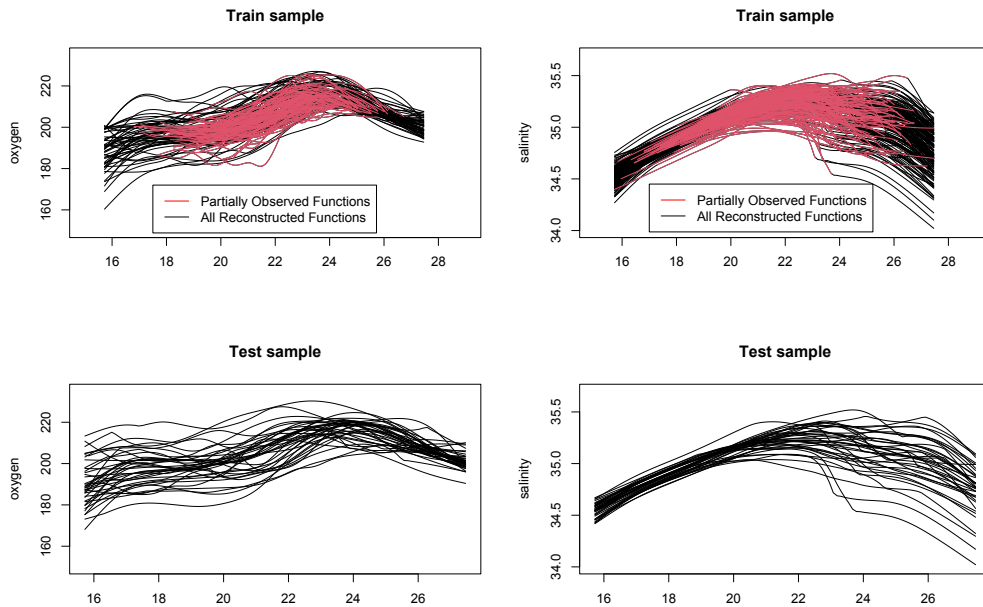


Figure 9: Training (resp. testing) samples of 153 (resp. 38) pairs of curves  $\{Oxygen_i, Salinity_i\}_{i=1}^{153(38)}$  in 'Scenario 2'.

Table 7: Real dataset: prediction errors over one hazard sample. MSPE, MSPA and  $k_n^*$  for 'Scenario 1' and 'Scenario 2'.

Methods		'SCENARIO 1'	'SCENARIO 2'
Rate of partially observed response curves (%)		12.042	52.356
Rate of missing data in response curve (%)		Less than 27	More than 45
<b>Reconst_X_Y</b>	MSPE	23.566	46.351
	MSPA	3.855	5.144
	$k_n^*$	13	8
<b>Reconst_X, Imp_Y</b>	MSPE	24.338	54.199
	MSPA	3.893	5.839
	$k_n^*$	9	12

## 7 Proofs

### 7.1 Proof of Theorem 3.1

Starting with the reconstruction cross covariance operator,

$$\begin{aligned}
\hat{\Delta}_{n,rec}^* &= \frac{1}{n} \sum_{i=1}^n Y_i^* \otimes X_i^* \\
&= \frac{1}{n} \sum_{i=1}^n \left( Y_i + (Y_i^* - Y_i) \right) \otimes X_i^*, \\
&= \frac{1}{n} \sum_{i=1}^n Y_i \otimes X_i^* + \frac{1}{n} \sum_{i=1}^n (Y_i^* - Y_i) \otimes X_i^*, \\
&= \Theta \hat{\Gamma}_{n,rec} + \frac{1}{n} \sum_{i=1}^n \epsilon_i \otimes X_i^* + \frac{1}{n} \sum_{i=1}^n (Y_i^* - Y_i) \otimes X_i^*.
\end{aligned}$$

Next, we obtain

$$\begin{aligned}
& \mathbb{E} \left( \left\| \hat{\Theta}^* \cdot X_{new}^* - \Theta \cdot X_{new}^* \right\|^2 \right) \\
&= \mathbb{E} \left( \left\| \hat{\Pi}_{k_n, rec} \hat{\Delta}_{n, rec}^* \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* - \Theta \cdot X_{new}^* \right\|^2 \right) \\
&\leq 4\mathbb{E} \left( \left\| \hat{\Pi}_{k_n, rec} \Theta \hat{\Gamma}_{n, rec} \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* - \Theta \cdot X_{new}^* \right\|^2 \right) \\
&+ 4\mathbb{E} \left( \left\| \hat{\Pi}_{k_n, rec} \left( \frac{1}{n} \sum_{i=1}^n \epsilon_i \otimes X_i^* \right) \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \right\|^2 \right) \\
&+ 2\mathbb{E} \left( \left\| \hat{\Pi}_{k_n, rec} \left( \frac{1}{n} \sum_{i=1}^n (Y_i^* - Y_i) \otimes X_i^* \right) \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \right\|^2 \right).
\end{aligned}$$

Applying several times the identity  $(a + b)^2 \leq 2a^2 + 2b^2$  for any  $a, b \in \mathbb{R}$ , we get

$$\begin{aligned}
\mathbb{E} \left( \left\| \hat{\Theta}^* \cdot X_{new}^* - \Theta \cdot X_{new}^* \right\|^2 \right) &\leq 64\mathbb{E} \left( \left\| \Theta \hat{\Pi}_{k_n, rec} X_{new}^* - \Theta \hat{\Pi}_{k_n} X_{new}^* \right\|^2 \right) \\
&+ 64\mathbb{E} \left( \left\| \Theta \hat{\Pi}_{k_n} X_{new}^* - \Theta \hat{\Pi}_{k_n} X_{new} \right\|^2 \right) \\
&+ 32\mathbb{E} \left( \left\| \Theta \hat{\Pi}_{k_n} X_{new} - \Theta \Pi_{k_n} X_{new} \right\|^2 \right) \\
&+ 16\mathbb{E} \left( \left\| \Theta \Pi_{k_n} X_{new} - \Theta X_{new} \right\|^2 \right) \\
&+ 8\mathbb{E} \left( \left\| \Theta X_{new} - \Theta X_{new}^* \right\|^2 \right) \\
&+ 4\mathbb{E} \left( \left\| \frac{1}{n} \sum_{i=1}^n \langle X_i^*, \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \rangle \epsilon_i \right\|^2 \right) \\
&+ 2\mathbb{E} \left( \left\| \frac{1}{n} \sum_{i=1}^n \langle X_i^*, \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \rangle (Y_i^* - Y_i) \right\|^2 \right).
\end{aligned}$$

Results of terms in the above decomposition are in [Crambes et al. \(2023\)](#), exceptionally the last term, let be noted by

$$P_n = \frac{1}{n} \sum_{i=1}^n \langle X_i^*, \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \rangle (Y_i^* - Y_i)$$

. Hence, using the Cauchy-Schwarz inequality, we have

$$\mathbb{E}(\|P_n\|^2) \leq \sqrt{\mathbb{E} \left( \left\| \frac{1}{n} \sum_{i=1}^n \langle X_i^*, \left( \hat{\Pi}_{k_n, rec} \hat{\Gamma}_{n, rec} \hat{\Pi}_{k_n, rec} \right)^{-1} X_{new}^* \rangle \right\|^4 \right)} \mathbb{E}(\|Y_i^* - Y_i\|^4).$$

The result comes from Lemma 5.2 in [Crambes and Henchiri \(2019\)](#) and the result (3.4) that gives us

$$\begin{aligned}\mathbb{E}(\|P_n\|^2) &= \mathcal{O}\left(\frac{k_n}{n}\right) + \mathcal{O}\left(n^{-\zeta_1(b_O-1)/(b_O+2)}\right) \\ &= \mathcal{O}\left(n^{\eta_1/(a_O+2)-1-\zeta_1(b_O-1)/(b_O+2)}\right).\end{aligned}$$

Summarizing, we get

$$\mathbb{E}\left(\left\|\hat{\Theta}^* \cdot X_{new}^* - \Theta \cdot X_{new}^*\right\|^2\right) = \mathcal{O}\left(n^{-\eta_1(a_O-1)/(2(a_O+2))} + n^{\eta_1/(a_O+2)-1-\zeta_1(b_O-1)/(b_O+2)}\right).$$

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