

Variable selection for functional linear models with strong heredity constraint

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Received: 15 August 2020 / Revised: 20 March 2021 / Accepted: 30 March 2021 /

Published online: 28 April 2021

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Abstract

In this paper, we consider the variable selection problem in functional linear regression with interactions. Our goal is to identify relevant main effects and corresponding interactions associated with the response variable. Heredity is a natural assumption in many statistical models involving two-way or higher-order interactions. Inspired by this, we propose an adaptive group Lasso method for the multiple functional linear model that adaptively selects important single functional predictors and pairwise interactions while obeying the strong heredity constraint. The proposed method is based on the functional principal components analysis with two adaptive group penalties, one for main effects and one for interaction effects. With appropriate selection of the tuning parameters, the rates of convergence of the proposed estimators and the consistency of the variable selection procedure are established. Simulation studies demonstrate the performance of the proposed procedure and a real example is analyzed to illustrate its practical usage.

Keywords Functional linear model · Main effect · Multiple functional predictors · Interaction effect · Heredity structure · Variable selection

1 Introduction

Functional data analysis has received considerable attentions in various fields, including, for example, environmental science, biology, medicine, finance and system engineering. The basic idea behind functional data analysis is to express each individual in repeatedly measured data as a smooth function and then make statistical inference from the collection of functional data (Ramsay & Silverman, 2005; Horváth & Kokoszka, 2012). There are a large body of recent work devoted to

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regression models with functional predictors, and among them, the most popular model is the functional linear model (Cardot et al., 2003; Cai & Hall, 2006; Hall & Horowitz, 2007; Crambes et al., 2009; Hall & Hooker, 2016). A standard functional linear model (FLM) with scalar response and functional predictor is defined as

$$Y = \alpha_0 + \int \beta(t)X(t)dt + \varepsilon, \tag{1}$$

where *Y* is the scalar response variable, X(t) is the random function defined on compact interval \mathcal{T} , α_0 is the intercept, $\beta(t)$ is the unknown coefficient function, and ε is the random error.

FLM is designed to describe the relation between a scalar response and one functional explanatory variable, which is often simple and easy to interpret. However, as pointed out in Yao and Müller (2010), this model imposes a constraint on the regression relationship that may not be appropriate in many real problems. In practice, the scalar response can be potentially associated with multiple or even a large number of functional predictors, which yields the multiple FLM as

$$Y = \alpha_0 + \sum_{j=1}^p \int \beta_j(t) X_j(t) dt + \varepsilon.$$
 (2)

For model (2), Matsui and Konishi (2011), Gertheiss et al. (2013) and Lian 2013 considered the shrinkage estimation and selection; Huang et al. (2016) studied the robust variable selection; Collazos et al. (2016) investigated the connection between model testing and variable selection; and Xue and Yao (2021) considered the hypothesis testing problem with the number of functional predictors being diverge as the sample size increases. In addition, (Kong et al., 2016; Ma et al., 2019) and (Yu et al., 2019) studied the regularized estimation and variable selection for the partial FLM with multiple functional predictors.

One limitation of model (2) is that the effects of the functional predictors are additive, that is, only the main effects of the individual functional predictors enter the regression model. In practice, however, ignoring the interaction effects may result in inaccurate or biased estimates of the model parameters which in turn lead to incorrect conclusions. To capture the interaction effects which can also arise in practice, it is often desired to develop new functional regression models that can accommodate both multiple functional predictors and the interaction terms among them. For this purpose, Yao and Müller (2010) considered the full quadratic effect $\int \int \gamma(s,t)X(s)X(t)dsdt$ where X(s) and X(t) represent the same functional predictor. Fuchs et al. (2015) and Usset et al. (2016) considered the interaction terms of the form $\int \int \gamma(s,t)X_1(s)X_2(t)dsdt$ where $X_1(s)$ and $X_2(t)$ represent different functional predictors.

In this paper, we consider the following functional regression model with main terms and all possible two-way interaction terms



$$Y = \alpha_0 + \sum_{j=1}^p \int \beta_j(t) X_j(t) dt + \sum_{j < m} \int \int \gamma_{jm}(s, t) X_j(s) X_m(t) ds dt + \varepsilon,$$
 (3)

where $X_j(t)$ $(j = 1, \dots, p)$ are main effects, and $X_j(s)X_m(t)$ $(1 \le j < m \le p)$ are twoway interaction effects. The regression coefficient functions $\beta_j(t)$ and $\gamma_{jm}(s,t)$ are assumed to be smooth and square integrable. When p = 2, model (3) reduces to the model in Fuchs et al. (2015) and Usset et al. (2016).

A special feature about model (3) is the intrinsic relationship among the regressor terms, for example, $X_i(s)X_m(t)$ is the child of $X_i(s)$ and $X_m(t)$, or equivalently, $X_i(s)$ and $X_m(t)$ are the parents of $X_i(s)X_m(t)$. This model structure is known as the hierarchical structure, with the two different types including strong heredity and weak heredity (Chipman, 1996; Bien et al., 2013). Strong heredity means that if an interaction term exists in the model, then both of its parent effects must be present; while for weak heredity, it only requires one of its parent effects must be present in the model. Heredity is a nature requirement in interaction models (Bien et al., 2013; Hao et al., 2018; She et al., 2018). There are two reasons one would prefer to add main effects to the model ahead of interaction effects. First, as pointed out by Cox (1984), large component main effects are more likely to lead to appreciable interactions than small components. And moreover, the interactions associated with larger main effects may also be in some sense of more practical importance. Second, interaction effects are often more difficult to interpret than main effects, thus, given similar predictive ability, one would prefer to add a main effect ahead of an interaction effect.

Variable selection for linear regression models with interaction effects has attracted considerable attentions in the past two decades. To name a few, Yuan et al. (2009) proposed non-negative garrote methods that naturally incorporate a general hierarchical structure among the predictors; Choi et al. 2010 extended the Lasso method for simultaneously fitting the regression model and identifying interaction terms under the strong heredity constraint; Bien et al. (2013) investigated a Lasso for hierarchical interactions; She et al. (2018) studied group regularized estimation under weak or strong heredity, respectively; Hao et al. (2018) proposed two-stage regularization methods on model selection and estimation for quadratic regression models under strong or weak heredity constraint, respectively. To the best of our knowledge, none of these papers studied the setting with functional data, and none of the variable selection methods for multiple functional linear models satisfied the heredity constraint when the interactions are also included.

In this paper, we propose to fill the gap by studying the variable selection and estimation of model (3) with strong heredity constraint. Based on the functional principal components analysis (FPCA) and the adaptive group Lasso (Yuan and Lin 2006), we propose a new variable selection method to adaptively select functional predictors and interactions while automatically obeying the strong heredity constraint. We carry out estimation and variable selection by optimizing a penalized least squares function that includes two adaptive penalty terms, one for the main effects and another for the reparameterized interaction effects. An effective algorithm has also been developed. With proper choice of tuning parameters, we



establish the convergence rates of the estimators for the coefficient functions and the consistency of the variable selection procedure.

The rest of the paper is organized as follows. In Sect. 2, we first describe the variable selection procedure using FPCA and the adaptive group Lasso penalties, and then propose an iterative algorithm for finding the penalized estimators. In Sect. 3, we present the theoretical properties of the proposed variable selection procedure. Simulation studies are carried out in Sect. 4 to assess the finite sample performance of the proposed estimators, and an environmental data set is analyzed in Sect. 5. Lastly, we conclude the paper in Sect. 6 with some future work. Technical details are given in the supplementary material.

2 Estimation methodology

2.1 Functional principal component analysis

Assume that we have independent and identically distributed observations $\{(Y_i, X_{ij}(t)), i = 1, \dots, n, j = 1, \dots, p\}$, where $X_{ij}(t)$ are zero mean random functions belonging to $L^2(T)$, and Y_i are response variables generated from model (3). We also assume that ε_i are independent and identically distributed random errors with a finite second moment, and they are independent of the functional predictors.

Let also $(X_1(t), \dots, X_p(t), Y)$ denote the generic random functions with the same distribution as $(X_{i1}(t), \dots, X_{ip}(t), Y_i)$ and denote the covariance function of $X_j(t)$ by $\Sigma_j(s,t) = \text{Cov}(X_j(s), X_j(t))$. Then by Mercer's Theorem, we can obtain the spectral decomposition as

$$\Sigma_{j}(s,t) = \sum_{k=1}^{\infty} \tau_{jk} \phi_{jk}(s) \phi_{jk}(t),$$

where $\tau_{j1} > \tau_{j2} > \cdots > 0$ are the eigenvalues of the linear operator associated with $\Sigma_j(s,t)$, and ϕ_{jk} are the corresponding eigenfunctions. By the Karhunen-Loève expansion, $X_j(t)$ can be represented as

$$X_j(t) = \sum_{k=1}^{\infty} \xi_{jk} \phi_{jk}(t),$$

where ξ_{jk} are the principal component scores satisfying $E(\xi_{jk}) = 0$, $E(\xi_{jk}^2) = \tau_{jk}$ and $E(\xi_{ik}\xi_{ik'}) = 0$ for $k \neq k'$. In addition, the sample $X_{ii}(t)$ can be expressed as

$$X_{ij}(t) = \sum_{k=1}^{\infty} \xi_{ijk} \phi_{jk}(t),$$

where ξ_{ijk} are independent copies of ξ_{jk} .

Since the sequences ϕ_{jk} $(k=1,2,\cdots)$ are all complete in the class of square integrable functions on \mathcal{T} , the regression coefficient functions in (3) can be represented as



$$\beta_{j}(t) = \sum_{k=1}^{\infty} \beta_{jk} \phi_{jk}(t), \quad \gamma_{jm}(s,t) = \sum_{k=1}^{\infty} \sum_{l=1}^{\infty} \gamma_{jm,kl} \phi_{jk}(s) \phi_{ml}(t), \tag{4}$$

for suitable sequences $(\beta_{jk})_{k=1,2,...}$ and $(\gamma_{jm,kl})_{k,l=1,2,...}$ with $\sum_k \beta_{jk} < \infty$ and $\sum_{k,l} \gamma_{jm,kl} < \infty$. By (4) and the orthonormality property of the eigenfunctions, model (3) can be alternatively expressed as a function of the scores ξ_{jk} and ξ_{ml} ,

$$Y_i = \alpha_0 + \sum_{j=1}^p \sum_{k=1}^\infty \beta_{jk} \xi_{ijk} + \sum_{j < m} \sum_{k=1}^\infty \sum_{l=1}^\infty \gamma_{jm,kl} \xi_{ijk} \xi_{iml} + \varepsilon_i.$$
 (5)

Our goals are to provide a method that determines which terms in the right-hand side of (5) have important effects on the response, and then construct the coefficient function estimators with desirable properties and develop an efficient computational algorithm. The existing variable selection methods for the functional regression models do not guarantee the strong heredity constraint, as they treat all the elements of $(\beta_{jk}, \gamma_{jm,kl})$ equally and do not distinguish between them. To tackle the problem, we first reparameterize the coefficients for the interaction terms $\gamma_{jm,kl}$ in (5) as $\gamma_{jm,kl} = \alpha_{jm,kl} \beta_{jk} \beta_{ml}$, which yields the strong hierarchical multiple functional linear model as

$$Y_i = \alpha_0 + \sum_{j=1}^p \sum_{k=1}^\infty \beta_{jk} \xi_{ijk} + \sum_{j < m} \sum_{k=1}^\infty \sum_{l=1}^\infty \alpha_{jm,kl} \beta_{jk} \beta_{ml} \xi_{ijk} \xi_{iml} + \varepsilon_i.$$
 (6)

With this reparameterization, the coefficient function for an interaction term $X_j(s)X_m(t)$ must be zero if either of its two main effects $X_j(s)$ or $X_m(t)$ has a zero coefficient function; and in contrast, if the coefficient function for $X_j(s)X_m(t)$ is not equal to zero, it implies that both $\beta_j(t) \neq 0$ and $\beta_m(t) \neq 0$, which guarantees the strong heredity constraint.

Note that the eigenvalue τ_{jk} of $X_j(t)$ often decreases to zero rapidly as k increases, that is, $X_j(t)$ mainly depends on the leading K_j principal components. This implies that it is reasonable to assume Y is dependent on the leading K_j principal components in $X_j(t)$. A practical strategy to select the smoothing parameters K_j will be discussed in Sect. 2.3. Moreover, for the sake of descriptive convenience, we set $\alpha_0 = 0$ in model (6) so that the final model is approximated as

$$Y_{i} \approx \sum_{j=1}^{p} \sum_{k=1}^{K_{j}} \beta_{jk} \xi_{ijk} + \sum_{j < m} \sum_{k=1}^{K_{j}} \sum_{l=1}^{K_{m}} \alpha_{jm,kl} \beta_{jk} \beta_{ml} \xi_{ijk} \xi_{iml} + \varepsilon_{i}.$$
 (7)

2.2 Model estimation

Since the scores ξ_{ijk} are unknown, we cannot estimate β_{jk} and α_{jm} directly. To estimate the functional principal component scores, we first estimate the covariance functions by



$$\hat{\Sigma}_{j}(s,t) = \frac{1}{n} \sum_{i=1}^{n} (X_{ij}(s) - \bar{X}_{j}(s))(X_{ij}(t) - \bar{X}_{j}(t)),$$

where $\bar{X}_{j}(s) = \frac{1}{n} \sum_{i=1}^{n} X_{ij}(s)$. Also by the empirical spectral expansion,

$$\hat{\Sigma}_{j}(s,t) = \sum_{k=1}^{\infty} \hat{\tau}_{jk} \hat{\phi}_{jk}(s) \hat{\phi}_{jk}(t),$$

where $\hat{\tau}_{j1} \geq \hat{\tau}_{j2} \geq \cdots \geq 0$ and $(\hat{\tau}_{jk}, \hat{\phi}_{jk})$ are pairs of eigenvalue and eigenfunction of $\hat{\Sigma}_{i}(s,t)$. Then consequently, the principal component scores ξ_{iik} can be estimated by

$$\hat{\xi}_{ijk} = \int (X_{ij}(s) - \bar{X}_j(s))\hat{\phi}_{jk}(s)\mathrm{d}s.$$

With these preliminary estimates at hand, we further denote

$$\begin{split} \hat{U}_{i} &= (\hat{U}_{i1}^{\top}, \cdots, \hat{U}_{ip}^{\top})^{\top}, \ \hat{U}_{ij} = (\hat{\xi}_{ij1}, \cdots, \hat{\xi}_{ijK_{j}})^{\top}, \ \hat{W}_{i} = (\hat{W}_{i12}^{\top}, \cdots, \hat{W}_{i(p-1)p}^{\top})^{\top}, \\ \hat{W}_{ijm} &= (\beta_{j1}\beta_{m1}\hat{\xi}_{ij1}\hat{\xi}_{im1}, \cdots, \beta_{j1}\beta_{mK_{m}}\hat{\xi}_{ij1}\hat{\xi}_{imK_{m}}, \cdots, \beta_{jK_{j}}\beta_{mK_{m}}\hat{\xi}_{ijK_{j}}\hat{\xi}_{imK_{m}})^{\top}, \\ \text{and } \beta &= (\beta_{1}^{\top}, \cdots, \beta_{p}^{\top})^{\top}, \beta_{j} = (\beta_{j1}, \cdots, \beta_{jK_{j}})^{\top}, \alpha = (\alpha_{12}^{\top}, \cdots, \alpha_{(p-1)p}^{\top})^{\top}, \\ \alpha_{jm} &= (\alpha_{jm,11}, \cdots, \alpha_{jm,1K_{m}}, \cdots, \alpha_{jm,K_{i}K_{m}})^{\top}, \end{split}$$

where $1 \le j < m \le p$. Then for the purpose of variable selection, we define the penalized least squares function with the adaptive group Lasso as

$$L(\beta, \alpha) = \sum_{i=1}^{n} (Y_i - \hat{U}_i^{\mathsf{T}} \beta - \hat{W}_i^{\mathsf{T}} \alpha)^2 + n \sum_{i=1}^{p} \lambda_{1j} \|\beta_j\|_2 + n \sum_{i < m} \lambda_{2,jm} \|\alpha_{jm}\|_2,$$
(8)

where $\|\cdot\|_2$ stands for the vector L_2 -norm. As can be seen, the tuning parameters λ_{1j} control the functional main effect estimates. If $\|\beta_j\|_2$ is shrunken to zero, all terms involving $X_j(t)$, including the main effect $\int \beta_j(t)X_j(t)dt$ and the interactions $\int \int \gamma_{jm}(s,t)X_j(s)X_m(t)dsdt$ for any m>j, will be removed from the model. The tuning parameters $\lambda_{2,jm}$ control the functional interaction effect estimates. If $\|\beta_j\|_2 \neq 0$ and $\|\beta_m\|_2 \neq 0$ but the corresponding interaction effect is not strong, $\|\alpha_{jm}\|_2$ still has the possibility of being zero. The penalty term controlled by $\lambda_{2,jm}$ thus provides the flexibility of selecting only main effects of $X_j(t)$ and $X_m(t)$ but not their interaction term.

2.3 Tuning parameter selection

For practical implementation, one has to decide the values of the tuning parameters and smoothing parameters. Note that there are a total of p(p + 1)/2 tuning parameters and p smoothing parameters in model (8), and thus the classical method including, for example, CV, AIC and BIC, may not be applicable.



To overcome this problem, we apply the methods in Zou (2006) and Wang et al. (2007) and simplify the tuning parameters as

$$\lambda_{1j} = \frac{\lambda}{\|\hat{\beta}_j^u\|_2^v} \quad \text{and} \quad \lambda_{2,jm} = \frac{\lambda}{\|\hat{\alpha}_{jm}^u\|_2^v},$$

where $\hat{\beta}^u_j$ and $\hat{\alpha}^u_{jm}$ are the unpenalized least squares estimators of β_j and α_{jm} , and v is a pre-specified positive number for which we take v=1 in our simulation study and real data analysis. Let also $\mathbf{K}=(K_1,\cdots,K_p)$. We then consider to select \mathbf{K} and the tuning parameter λ according to the following extended BIC (EBIC) criterion:

$$EBIC(\lambda, \mathbf{K}) = \log \left\{ \frac{1}{n} \sum_{i=1}^{n} (Y_i - \hat{U}_i^{\mathsf{T}} \hat{\beta} - \hat{W}_i^{\mathsf{T}} \hat{\alpha})^2 \right\} + df \frac{\log n}{n} + 2\kappa \frac{\log P}{n}, \quad (9)$$

where P is the total dimension of the model space, and $\kappa \in [0, 1]$ is a regulation parameter, and the associated degree of freedom is

$$\begin{split} df &= \sum_{j=1}^{p} I \Big\{ \| \hat{\beta}_{j} \|_{2} > 0 \Big\} + \sum_{j=1}^{p} \frac{\| \hat{\beta}_{j} \|_{2}}{\| \hat{\beta}_{j}^{u} \|_{2}} (K_{j} - 1) \\ &+ \sum_{j < m} I \Big\{ \| \hat{\alpha}_{jm} \|_{2} > 0 \Big\} + \sum_{j < m} \frac{\| \hat{\alpha}_{jm} \|_{2}}{\| \hat{\alpha}_{jm}^{u} \|_{2}} (K_{j} K_{m} - 1). \end{split}$$

According to Chen and Chen (2008), we take $\kappa = 0.5$ throughout the simulation study and real data analysis. To further reduce the computations of $\mathrm{EBIC}(\lambda, \mathbf{K})$, one can set $K_j \equiv K_0$ for all $1 \le j \le p$ so that the resulting criterion reduces to $\mathrm{EBIC}(\lambda, K_0)$.

Another way to alleviate the computation burden is to apply a two-stage method that first selects K_j by the cumulative percentage of total variance (CPTV) criterion and then chooses λ by EBIC. To be more specific, we first define the CPTV explained by the first d functional principal components as

$$CPTV_{j}(d) = \frac{\sum_{l=1}^{d} \hat{\tau}_{jl}}{\sum_{l=1}^{n} \hat{\tau}_{il}},$$
(10)

and choose K_j as the minimum number of d for which $CPTV_j(d)$ exceeds a desired level $100\delta_0\%$. In our numerical studies, we adopt $\delta_0=0.95$. Denote also the selected smoothing parameters by CPTV as $\mathbf{K}_{CPTV}=(K_1,\cdots,K_p)$. Then as the second step, we select λ by the EBIC criterion in (9) with \mathbf{K}_{CPTV} given. For convenience, we refer to this selection method as the C-EBIC criterion.

Despite the simplified tuning parameter selection may not be optimal, the resulting estimate of λ does guarantee that the tuning parameter for zero coefficient is larger than that for nonzero coefficient. Thus, we can consistently estimate the large coefficients and shrink the small coefficients toward zero simultaneously. From the numerical studies in Sect. 4, we can see that the proposal tuning parameter selection performs very well in practice.



2.4 Computational algorithm

Because the main effect coefficients β and the interaction coefficients α are controlled at different levels, we apply an iterative algorithm to estimate them alternatively. Specifically, we first fix β_j and estimate α_{jm} , then fix α_{jm} and estimate β_j , and finally iterate between these steps until convergence. When β is fixed, the optimization in α becomes a group Lasso problem, hence one can use either the group LARS algorithm (Yuan and Lin, 2006) or the quadratic programming to solve for α efficiently. When α is fixed, we solve for β_1, \dots, β_p sequentially. For each $j = 1, \dots, p$, we fix α and $\beta_{[j]} = (\beta_1^\top, \dots, \beta_{j-1}^\top, \beta_{j+1}^\top, \dots, \beta_p^\top)^\top$, then (8) becomes a simple group Lasso problem with only one coefficient β_i .

Specifically, for a fixed λ , the proposed iterative algorithm is as follows:

Step 1. Find the initial estimators $\hat{\beta}^{(0)}$ and $\hat{\alpha}^{(0)}$. For example, the unpenalized estimators obtained by minimizing (8) with $\lambda = 0$ can be used.

Step 2. Fixing $\hat{\beta}^{(0)}$, we update $\hat{\alpha}$ by

$$\hat{\alpha}^{(1)} = \arg\min_{\alpha} \left\{ \sum_{i=1}^{n} (Y_i - \hat{U}_i^{\mathsf{T}} \hat{\beta}^{(0)} - \hat{W}_i^{(0)\mathsf{T}} \alpha)^2 + n\lambda \sum_{i \le m} \frac{\|\alpha_{jm}\|_2}{\|\hat{\alpha}_{im}^u\|_2} \right\},\tag{11}$$

where $\hat{W}_{i}^{(0)}$ has the same form as \hat{W}_{i} except that $\beta_{jk}\beta_{ml}$ are replaced by $\hat{\beta}_{jk}^{(0)}\hat{\beta}_{ml}^{(0)}$. Step 3. Fixing $\hat{\alpha}^{(1)}$, we update $\hat{\beta}_{a}$ by

$$\hat{\beta}_{q}^{(1)} = \arg\min_{\beta_{q}} \left\{ \sum_{i=1}^{n} (\tilde{Y}_{i} - \sum_{k=1}^{K_{q}} \beta_{qk} \tilde{U}_{iqk})^{2} + n\lambda \frac{\|\beta_{q}\|_{2}}{\|\hat{\beta}_{q}^{u}\|_{2}} \right\}, \tag{12}$$

where $q = 1, 2, \dots, p$ and

$$\begin{split} \tilde{Y}_i &= Y_i - \sum_{j=1, j \neq q}^p \sum_{k=1}^{K_j} \hat{\beta}_{jk}^{(0)} \hat{\xi}_{ijk} - \sum_{j < m, j, m \neq q} \sum_{k=1}^{K_j} \sum_{l=1}^{K_m} \hat{\alpha}_{jm, kl}^{(1)} \hat{\beta}_{jk}^{(0)} \hat{\beta}_{ml}^{(0)} \hat{\xi}_{ijk} \hat{\xi}_{iml}, \\ \tilde{U}_{iqk} &= \hat{\xi}_{iqk} + \sum_{l=1}^{q-1} \sum_{k=1}^{K_j} \hat{\alpha}_{jq, hk}^{(1)} \hat{\beta}_{jh}^{(0)} \hat{\xi}_{ijh} \hat{\xi}_{iqk} + \sum_{l=1}^p \sum_{k=1}^{K_m} \hat{\alpha}_{qm, kh}^{(1)} \hat{\beta}_{mh}^{(0)} \hat{\xi}_{iqk} \hat{\xi}_{imh}. \end{split}$$

Step 4. Setting $\hat{\beta}^{(0)} = \hat{\beta}^{(1)}$ and $\hat{\alpha}^{(0)} = \hat{\alpha}^{(1)}$, we iterate Step 2 and Step 3 until convergence to obtain the final estimator $\hat{\theta} = (\hat{\beta}^{\mathsf{T}}, \hat{\alpha}^{\mathsf{T}})^{\mathsf{T}}$, where the convergence criterion is set as $\|\hat{\theta}^{(1)} - \hat{\theta}^{(0)}\|_2 < 10^{-4}$.

We minimize the objective function $L(\beta,\alpha)$ with respect to either the set of α 's or the set of β 's so that the objective function decreases at each step. The value of the objective function is then guaranteed to converge to a local minimum since it is bounded from below. Similar to many penalized estimation problems, a convergence of the algorithm to the global minimum may not be guaranteed, although the simulation results did show that our proposed algorithm can effectively and accurately detect the true model structure.

Lastly, for each $j, m = 1, 2, \dots, p$ with j < m, we estimate $\beta_i(t)$ and $\gamma_{im}(s, t)$ by



$$\hat{\beta}_{j}(t) = \sum_{k=1}^{K_{j}} \hat{\beta}_{jk} \hat{\phi}_{jk}(t) \quad \text{and} \quad \hat{\gamma}_{jm}(s,t) = \sum_{k=1}^{K_{j}} \sum_{l=1}^{K_{m}} \hat{\alpha}_{jm} \hat{\beta}_{jk} \hat{\beta}_{ml} \hat{\phi}_{jk}(s) \hat{\phi}_{ml}(t).$$

3 Asymptotic properties

We assume that the true model obeys the strong heredity constraint, and let β_j^* and α_{jm}^* denote the underlying true values of β_j and α_{jm} . It is noted that $\beta_j^*(t)$ is zero if and only if $\beta_j^*=0$, and $\gamma_{jm}^*(s,t)$ is nonzero only if α_{jm}^* , β_j^* and β_m^* are all nonzero vectors. We further define

$$A_1 = \{j : \beta_j^*(t) \neq 0\}$$
 and $A_2 = \{(j, m) : \gamma_{jm}^*(s, t) \neq 0\},$

where A_1 contains the indices for main terms whose true coefficient functions are nonzero, and A_2 contains the indices for interaction terms whose true coefficient functions are nonzero. Let also

$$\begin{split} &a_n = \max\{\lambda_{1j}, \lambda_{2,mm'} \, : \, j \in A_1, (m,m') \in A_2\}, \\ &b_n = \min\{\lambda_{1j}, \lambda_{2,mm'} \, : \, j \in A_1^c, (m,m') \in A_2^c, m, m' \in A_1\}. \end{split}$$

In addition, for convenience and simplicity, we let $\|\cdot\|$ represent the $L^2(\mathcal{T})$ norm and C denote a positive constant that may be different at each appearance throughout this paper.

To derive the asymptotic properties of the variable selection procedure and the corresponding estimators, we need the following regularity conditions.

- (A1) $X_{ij}(s)$ has a finite fourth moment such that $\int E(X_{ij}^4(s)) ds < \infty$ for $j = 1, \dots, p$.
- (A2) There exist some constants C > 1, a > 1 and b > a + 1/2 such that for any $k \ge 1$,

$$C^{-1}k^{-a} \le \tau_{jk} \le Ck^{-a}, \quad \tau_{jk} - \tau_{j,k+1} \ge C^{-1}k^{-a-1} \quad \text{and} \quad |\beta_{jk}| \le Ck^{-b}.$$

- (A3) Assume that $E(\xi_{jk}^4) \le C\tau_{jk}^2$ for some constant C > 1, $E(\xi_{jk}^{v_1}\xi_{jl}^{v_2}) = E(\xi_{jk}^{v_1})E(\xi_{jl}^{v_2})$ for $v_1 + v_2 = 3$ and $1 \le k, l < \infty$, and $E(\xi_{jk}^{v_1}\xi_{jl}^{v_2}) = E(\xi_{jk}^{v_1})E(\xi_{jl}^{v_2})$ for $v_1 + v_2 = 4$ and $1 \le k \ne l < \infty$.
- (A4) $K_1 \times K_2 \cdots \times K_p \times K$, and $K = O(n^{1/(a+2b)})$, where $a_0 \times b_0$ means that the ratio a_0/b_0 is bounded away from zero and infinity.
- (A5) $\sqrt{nK^{-1}}a_n \to 0$ and $\sqrt{nK^{-1}}b_n \to \infty$, as $n \to \infty$.

Condition (A1) is an assumption on the moments, and condition (A2) is imposed upon the rate of the eigenvalues of the covariance operator of the functional predictors. Both (A1) and (A2) are commonly assumed in functional linear regression, see, for example, Hall and Horowitz (2007) and Cai and Hall (2006). Condition (A3) is similar to that of Yao and Müller (2010), condition (A4) is similar to



that of Hall and Horowitz (2007), and condition (A5) is similar to that of Choi et al. (2010). Note also that conditions (A4) and (A5) are needed for the consistency of the estimators and for the model detection.

Theorem 1 Assume that conditions (A1)–(A4) hold. Then

(i) if
$$\sqrt{K}a_n = o(1)$$
 as $n \to \infty$, we have
$$\|\hat{\beta}_j(t) - \beta_j^*(t)\|^2 = O_p(K(\sqrt{K/n} + a_n)^2 + K^{-2b+1});$$

(ii) if $Ka_n = o(1)$ as $n \to \infty$, we have $\|\hat{\gamma}_{jm}(s,t) - \gamma_{jm}^*(s,t)\|^2 = O_p(K^2(\sqrt{K/n} + a_n)^2 + K^{-4b+2}).$

The proof of Theorem 1 is given in the supplementary material. Theorem 1 demonstrates how the rates of convergence of the penalized estimators depend on λ_{1j} , $\lambda_{2,jm}$ and K. Moreover, if condition (A5) also holds, the rates of convergence in Theorem 1 will become

$$\|\hat{\beta}_j(t) - \beta_j^*(t)\|^2 = O_p(n^{-\frac{2b-1}{a+2b}}) \quad \text{and} \quad \|\hat{\gamma}_{jm}(s,t) - \gamma_{jm}^*(s,t)\|^2 = O_p(n^{-\frac{a+2b-3}{a+2b}}).$$

As in Hall and Horowitz (2007) for functional linear regression, the convergence rates of $\hat{\beta}_j(t)$ and $\hat{\gamma}_{jm}(s,t)$ are determined by the smoothness levels of the coefficient function and the covariance function, respectively.

Theorem 2 Assume that conditions (A1)–(A5) hold. As $n \to \infty$, we have

$$P(\hat{\beta}_j(t) = 0 \text{ for } j \in A_1^c) \to 1 \text{ and } P(\hat{\gamma}_{jm}(s, t) = 0 \text{ for } (j, m) \in A_2^c) \to 1.$$

The proof of Theorem 2 is given in the supplementary material. Theorem 2 shows that our proposed method can consistently remove the noise terms with probability tending to 1. That is, as long as the sample size is sufficiently large, our method is able to select the true model with a high probability.

4 Simulation studies

In this section, we conduct simulation studies to assess the finite sample performance of the proposed variable selection procedure. Specifically for $t \in [0, 1]$, we first generate the functional predictors from the following processes:



$$\begin{split} X_{i1}(t) &= \xi_{i11}\phi_1(t) + \xi_{i12}\phi_2(t), \\ X_{i2}(t) &= \xi_{i21}\phi_1(t) + \xi_{i22}\phi_2(t) + \xi_{i23}\phi_3(t), \\ X_{i3}(t) &= \xi_{i31}\phi_1(t) + \xi_{i32}\phi_4(t) + \xi_{i33}\phi_6(t), \\ X_{i4}(t) &= \xi_{i41}\phi_5(t) + \xi_{i42}\phi_6(t), \end{split}$$

where $\phi_k(t) = \sqrt{2}\cos(0.5(k+1)\pi t)$ for k = 1, 3 and $5, \phi_k(t) = \sqrt{2}\sin(0.5k\pi t)$ for k = 2, 4 and 6, and the random coefficients are as follows:

$$\begin{split} \xi_{i1l} &\sim N(0,\tau_{1l}) \text{ with } \tau_{11} = 4, \tau_{12} = 1/2, \\ \xi_{i2l} &\sim N(0,\tau_{2l}) \text{ with } \tau_{21} = 4, \tau_{22} = 1, \tau_{23} = 1/2, \\ \xi_{i3l} &\sim N(0,\tau_{3l}) \text{ with } \tau_{31} = 4, \tau_{32} = 2, \tau_{33} = 1, \\ \xi_{i4l} &\sim N(0,\tau_{4l}) \text{ with } \tau_{41} = 4, \tau_{42} = 1. \end{split}$$

We then generate the response variables Y_i from the model:

$$Y_{i} = \sum_{i=1}^{4} \int_{0}^{1} \beta_{j}(t) X_{ij}(t) dt + \sum_{i=1}^{3} \sum_{m=i+1}^{4} \int_{0}^{1} \int_{0}^{1} \gamma_{jm}(s, t) X_{ij}(s) X_{im}(t) ds dt + \varepsilon_{i},$$

where $\beta_1(t) = -0.5\phi_1(t) + 2\phi_2(t)$, $\beta_2(t) = -0.75\phi_1(t) + 1.5\phi_2(t) + 0.5\phi_3(t)$, $\beta_3(t) = -0.25\phi_1(t) + 0.5\phi_2(t) + 0.5\phi_3(t) + 0.5\phi_4(t)$, $\beta_4(t) = 0$, and ε_i are independent errors from N(0, 0.25). Finally, we consider two models in our simulations, one with interaction effects and one without interaction effect, as follows:

Model (I):
$$\gamma_{12}(s,t) = \frac{1}{3}\phi_2(s)\phi_1(t) + \frac{1}{\sqrt{3}}\phi_4(s)\phi_3(t)$$
, and $\gamma_{13}(s,t) = \gamma_{14}(s,t) = \gamma_{23}(s,t) = \gamma_{24}(s,t) = \gamma_{34}(s,t) = 0$;

Model (II):
$$\gamma_{12}(s,t) = \gamma_{13}(s,t) = \gamma_{14}(s,t) = \gamma_{23}(s,t) = \gamma_{24}(s,t) = \gamma_{34}(s,t) = 0$$
.

Following the above simulation settings, the random samples (Y_i, X_{ij}) are obtained with each X_{ij} being observed at 100 equally spaced points on [0, 1]. Moreover, we assume that the observations of the random trajectory X_{ij} at each point $t_{ij,r}$ are contaminated with measurement error $\epsilon_{ij,r}$ from N(0, 0.04). We compare the performance of our model selection procedure with the group SCAD (Lian, 2013) and the group Lasso (Gertheiss et al., 2013), which do not guarantee the strong heredity constraint. The sample sizes in our simulations are n=100, 200 and 300, respectively. Further for model selection, we apply both $\mathrm{EBIC}(\lambda, K_0)$ and C-EBIC in Sect. 2.3 to select the tuning parameter λ and the smoothing parameters K_i .

We carried out 500 simulations for each setting and then summarize the results in Tables 1 and 2 for model selection by the three methods under the two models, respectively. The column labeled " C_M " gives the average number of the nonzero main effects that are correctly selected, " C_I " gives the average number of the nonzero interactions that are correctly selected, and the column labeled " C_Z " gives the average number of the six true zeros that are correctly set to zero. The columns "UF", "CF" and "OF" are the proportions of models that are underfitted, correctly fitted and overfitted, respectively.

From Table 1, we can see that the proposed method correctly selects the true model more frequently than the group SCAD and the group Lasso. Specifically, for all scenarios, the group SCAD and the group Lasso perform similar, with the



Table 1 Variable selection results of the three methods under Model (I)

		Proposed method			Group SCAD			Group Lasso		
	n	C_M	C_I	C_Z	C_M	C_I	C_Z	C_M	C_I	C_Z
	100	2.9740	0.9660	5.9020	2.1160	0.8040	4.2080	1.9820	0.7740	3.8260
EBIC	200	2.9820	0.9880	5.9780	2.0320	0.8180	5.4540	1.9960	0.8120	5.1240
	300	3	1	5.9940	2.1020	0.8560	5.7480	2.0140	0.8420	5.5520
	100	2.9780	0.9780	5.9400	2.0280	0.7820	4.0880	1.9420	0.7620	3.9280
C-EBIC	200	2.9880	0.9900	5.9760	2.0380	0.8160	5.4760	1.9800	0.8020	5.1460
	300	3	1	5.9920	2.0600	0.8600	5.7540	2.0060	0.8340	5.5440
	n	UF	CF	OF	UF	CF	OF	UF	CF	OF
	100	0.0880	0.8240	0.0880	0.3620	0.2860	0.3520	0.3960	0.0480	0.5560
EBIC	200	0.0460	0.9120	0.0420	0.3660	0.3060	0.3280	0.3760	0.0560	0.5680
	300	0.0120	0.9780	0.0100	0.2480	0.3540	0.3980	0.2720	0.0520	0.6760
	100	0.0920	0.8260	0.0820	0.3540	0.2840	0.3620	0.3880	0.0360	0.5760
C-EBIC	200	0.0440	0.9060	0.0500	0.3640	0.3100	0.3260	0.3420	0.0420	0.6160
	300	0	0.9740	0.0260	0.2540	0.3540	0.3920	0.2380	0.0340	0.7280

Table 2 Variable selection results of the three methods under Model (II)

		Propose	Proposed method		Group SCAD			Group Lasso		
	n	C_M	C_I	C_Z	C_M	C_I	C_Z	C_M	C_I	C_Z
	100	2.9360	0.1280	6.7820	2.9780	0.0580	6.9780	2.9420	0.0860	6.8120
EBIC	200	2.9760	0.0320	6.9060	2.9920	0.0220	6.9860	2.9720	0.0560	6.9120
	300	2.9820	0.0240	6.9540	3	0	6.9920	3	0	6.9580
C-EBIC	100	2.9420	0.1320	6.7680	2.9880	0.0520	6.9800	2.9380	0.1020	6.7940
	200	2.9740	0.0480	6.9140	2.9940	0.0160	6.9880	2.9760	0.0540	6.8860
	300	2.9880	0.0160	6.9520	3	0	6.9960	2.9860	0	6.9620
	n	UF	CF	OF	UF	CF	OF	UF	CF	OF
	100	0.0440	0.8540	0.1020	0.0160	0.9180	0.0660	0.0440	0.8760	0.0800
EBIC	200	0.0240	0.9160	0.0600	0.0080	0.9640	0.0280	0.0260	0.9120	0.0620
	300	0.0160	0.9680	0.0160	0.0040	0.9940	0.0020	0.0180	0.9740	0.0080
	100	0.0420	0.8520	0.1060	0.0120	0.9200	0.0680	0.0420	0.8740	0.0840
C-EBIC	200	0.0220	0.9120	0.0660	0.0060	0.9620	0.0320	0.0180	0.9120	0.0700
	300	0.0060	0.9720	0.0220	0	0.9960	0.0040	0.0120	0.9720	0.0160

former slightly better, while our proposed method performs much better than both of them. When the sample size increases, the proposed method shows a consistent model selection trend. In addition, we note that $\mathrm{EBIC}(\lambda, K_0)$ and C-EBIC perform nearly the same. Lastly, Table 2 indicates that when there is no interaction effect (Model (II)), the proposed method performs also comparably to the other two methods.



To assess the performance of estimators $\hat{\beta}_1(\cdot)$, $\hat{\beta}_2(\cdot)$, $\hat{\beta}_3(\cdot)$ and $\hat{\gamma}_{12}(\cdot, \cdot)$, we apply the mean integrated squared errors (MISEs):

MISE_j =
$$\int_0^1 {\{\hat{\beta}_j(t) - \beta_j(t)\}^2 dt},$$

MISE₁₂ = $\int_0^1 \int_0^1 {\{\hat{\gamma}_{12}(t,s) - \gamma_{12}(t,s)\}^2 dt ds}.$

Table 3 presents the median of MISE for the coefficient functions $\hat{\beta}_j(\cdot)$ and $\hat{\gamma}_{12}(\cdot,\cdot)$ over the 500 simulations. The column labeled "Oracle" means the oracle estimators computed by the true model. From Table 3, we can see that all the estimators of the coefficient functions are close to the true curves. As the sample size n increases, the MISEs of all the estimators decrease. We also note that the performance of C-EBIC is slightly better than $\mathrm{EBIC}(\lambda, K_0)$. One possible explanation is that for $\mathrm{EBIC}(\lambda, K_0)$, the numbers of functional principal components for each predictor are restricted to be the same; while for C-EBIC, it allows different predictors to have different numbers of functional principal components.

5 Real data application

For illustration purpose, we apply the proposed method to the air pollution data of Beijing. The data are freely available on UCI Machine Learning Repository (with link https://archive.ics.uci.edu/ml/datasets/Beijing+Multi-Site+Air-Quality+Data). The data set consists of a collection of hourly measurements of air pollutants and

			Proposed method			Oracle			
			n = 100	n = 200	n = 300	n = 100	n = 200	n = 300	
Model (I)		MISE ₁	0.0171	0.0093	0.0095	0.0093	0.0058	0.0067	
		$MISE_2$	0.0399	0.0292	0.0341	0.0155	0.0097	0.0084	
	EBIC	$MISE_3$	0.1036	0.1102	0.1029	0.0998	0.0986	0.0974	
		$MISE_{12}$	0.2214	0.2324	0.2007	0.1363	0.1077	0.0953	
		$MISE_1$	0.0128	0.0063	0.0058	0.0074	0.0038	0.0026	
	C-EBIC	$MISE_2$	0.0338	0.0264	0.0235	0.0120	0.0064	0.0047	
		$MISE_3$	0.0960	0.0939	0.0925	0.0750	0.0727	0.0719	
		$MISE_{12}$	0.2188	0.2298	0.1991	0.1316	0.1045	0.0914	
Model (II)		$MISE_1$	0.0225	0.0158	0.0144	0.0095	0.0092	0.0069	
	EBIC	$MISE_2$	0.0416	0.0314	0.0291	0.0196	0.0188	0.0187	
		$MISE_3$	0.1037	0.0975	0.0923	0.0842	0.0686	0.0701	
		$MISE_1$	0.0202	0.0118	0.0113	0.0079	0.0068	0.0034	
	C-EBIC	$MISE_2$	0.0368	0.0297	0.0252	0.0187	0.0157	0.0143	
		MISE ₃	0.0973	0.0809	0.0776	0.0728	0.0580	0.0523	



weather factors in Beijing between January 1, 2015 and December 31, 2015, where the daily average PM2.5 concentration is of interest and set as the scalar response. The eight functional predictors, $X_1(t)$, $X_2(t)$, \cdots , $X_8(t)$, are, respectively, the hourly observed concentration of respirable suspended particulate (μ g/m³, PM10), sulphur dioxide (μ g/m³, SO₂), nitrogen dioxide (μ g/m³, NO₂), carbon monoxide (μ g/m³, CO), and ozone (μ g/m³, O₃), and the hourly observed meteorological variables including temperature (in Celsius, TEMP), dew point temperature (in Celsius, DEWP), and wind speed (m/s, WSPM).

In this section, our aim is to study the relationship of the daily average PM2.5 concentration to the air pollution and the weather conditions, and select the functional predictors and their interactions that contribute the most to the prediction of the daily average PM2.5 concentration. To be more specific, we propose the following functional interaction model with 8 functional predictors and 28 interaction terms, which naturally capture the cumulative effects and the interaction effects:

$$Y = \sum_{j=1}^{8} \int \beta_j(t) X_j(t) dt + \sum_{l < m} \int \int \gamma_{lm}(s, t) X_l(s) X_m(t) ds dt + \varepsilon, \tag{13}$$

where *Y* is the logarithm of the daily average PM2.5 concentration. In order to evaluate the performance of the proposed method, we use the first 300 observations as training sample and the last 65 observations as test sample, where the training sample is used to select the significant variables and estimate the coefficient functions, and the test sample is used to verify the quality of prediction. Finally, we apply the following mean squared error of prediction (MSEP) as the criterion for comparison:

MSEP =
$$\frac{1}{65} \sum_{i=301}^{365} (Y_i - \hat{Y}_i)^2$$
.

Taking into account the computation efficiency of C-EBIC and its good performance in simulation studies as shown in Sect. 4, we adopt the C-EBIC criterion to select the smoothing parameters K_i and the tuning parameter λ . That is, we first select K_i by the CPTV criterion in (10), and then select the tuning parameter λ by the EBIC criterion in (9) with K_i given. The significant variables selected by the proposed method and the group Lasso are presented in Table 4. We note that the both methods selected the main effects PM10, SO₂, NO₂, O₃, TEMP and WSPM and the interaction effects PM10 \times WSPM, SO₂ \times WSPM and NO₂ \times WSPM. This shows that the formation of PM2.5 is affected by multi-factors. The group Lasso also selected interaction effects DEWP \times WSPM, DEWP \times PM10 and CO \times NO₂, but they do not obey the heredity constraint; on the other hand, our proposed method selected the main effect DEWP but the group Lasso did not. The MSEPs of the selected models and the main effect model are displayed in Table 5. Figure 1 presents the original data and predicted values of the test sample. According to Fig. 1, the proposed model and estimation method perform well in predicting the PM2.5 concentration.

To further assess the stability of model selection, we apply the bootstrap analysis on the air pollution data. Based on 300 bootstrap samples of the training sample,



Table 4 Selected main effects and interactions in the air pollution data

Proposed met	hod	Group Lasso			
Main effects	Interactions	Main effects	Interactions		
PM10	PM10×WSPM	PM10	PM10×WSPM		
SO_2	$SO_2 \times WSPM$	SO_2	$SO_2 \times WSPM$		
NO_2	$NO_2 \times WSPM$	NO_2	$NO_2 \times WSPM$		
O_3	$O_3 \times WSPM$	O_3	$DEWP \times WSPM$		
TEMP	$TEMP \times WSPM$	TEMP	$DEWP \times PM10$		
DEWP	$\text{DEWP} \times \text{WSPM}$	WSPM	$CO \times NO_2$		
WSPM					

Table 5 The MSEPs for the air pollution data

	Main effect model	Interaction model			
		Proposed method	Group Lasso		
Training sample	0.2954	0.1621	0.1933		
Bootstrap sample	0.3437	0.2731	0.2964		
(standard deviations)	(0.0816)	(0.0752)	(0.0808)		

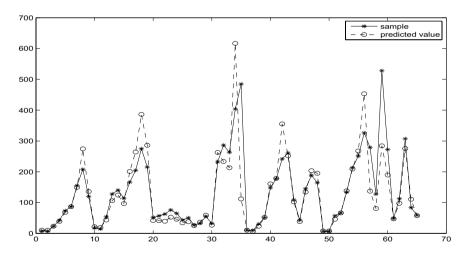


Fig. 1 The sample observations and predicted values

the main effects and interactions with selection frequency (denoted briefly as "Fre") higher than 30% are summarized in Table 6. These results indicate that the proposed method is fairly stable in terms of selecting terms. The average MSEP (and standard deviation) of the bootstrap samples are also displayed in Table 5, and from which it is evident that the performance of the proposed method is better than the group Lasso, and meanwhile both of them perform much better than the main effect model.



Proposed method				Group Lasso				
Main effects	Fre (%)	Interactions	Fre (%)	Main effects	Fre (%)	Interactions	Fre (%)	
PM10	100	PM10×WSPM	92	PM10	100	PM10×WSPM	90	
SO_2	62	$SO_2 \times WSPM$	51	SO_2	61	$SO_2 \times WSPM$	37	
NO_2	98	$NO_2 \times WSPM$	79	NO_2	83	$NO_2 \times WSPM$	65	
O_3	92	$O_3 \times WSPM$	86	O_3	85	$DEWP \times WSPM$	54	
TEMP	48	TEMP× WSPM	37	TEMP	38	$DEWP \times PM10$	37	
DEWP	37	$\text{DEWP} \times \text{WSPM}$	31	WSPM	75	$CO \times NO_2$	33	
WSPM	95	$PM10 \times DEWP$	36	DEWP	33	$NO_2 \times O_3$	35	
						$\mathrm{PM10}\times\mathrm{NO}_2$	38	

Table 6 Variable selection results based on the bootstrap samples

To conclude, by using the proposed variable selection procedure with the inclusion of possible interaction effects, we can obtain a more interpretable model with a reasonably good prediction performance.

Finally, we use all the 365 observations to conduct the variable selection. The final model contains 7 main effects: PM10, SO₂, NO₂, O₃, TEMP, DEWP and WSPM, and 6 interaction effects: PM10 × WSPM, SO₂ × WSPM, NO₂ × WSPM, O₃ × WSPM, TEMP × WSPM and DEWP × WSPM. Also for visualization, we display the estimated coefficient functions for the main effects in Fig. 2, and the estimated coefficient functions for the interaction effects in Fig. 3. From the figures, we have a few interesting findings as follows. First, the gaseous pollutants PM10, NO₂, O₃ and

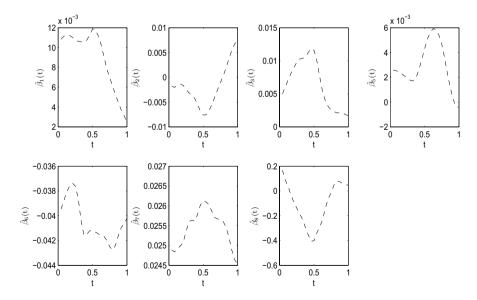


Fig. 2 Estimated coefficient functions for the main effects



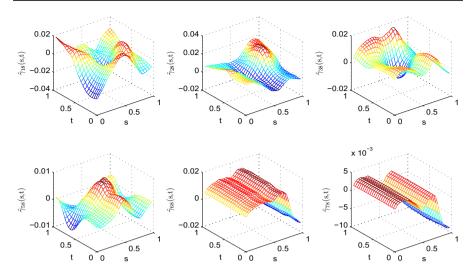


Fig. 3 Estimated coefficient functions for the interaction effects

the meteorological variable DEWP have a positive effect on the PM2.5 level. This conclusion is similar to the result in Wan et al. (2021). Second, TEMP has a significantly negative effect on the PM2.5 level. This finding is consistent with the chemical transport model sensitivity study in Dawson et al. (2007). Third, WSPM displays a more complex relationship with PM2.5, which may depend on other factors such as the wind direction (Tai et al., 2010; Wan et al. 2021).

6 Conclusion

In this paper, we extended the adaptive group Lasso method to accommodate the multiple functional linear models including the interaction terms. The proposed method automatically obeys the strong heredity constraint. With appropriate selection of the regularization parameters, we established the convergence rates of the estimators for the coefficient functions and the consistency of the variable selection procedure. Simulation studies indicated that our new method is able to consistently select the true model and works efficiently in estimation with finite samples. A real data example was also analyzed to illustrate the usefulness of our method in practice.

There are some interesting future directions. In this paper, we only considered the strong heredity, that is, an interaction is allowed only if both of the associated main effects are present in the model. As mentioned in Sect. 1, however, another common heredity can be the weak heredity for which only one of the main effects is required to be present in the model for their interaction to be included. As a future work, we will extend our new procedure to obey the weak heredity constraint and investigate its statistical properties and applications. Another interesting direction can be to extend our new procedure to the generalized functional linear models with strong or weak heredity constraint.



7 Supplementary Material

Supplement to "Variable selection for functional linear models with strong heredity constraint". To save space, the proofs of Theorems 1–2, Lemmas 1–2 and their proofs, and additional simulation results are provided in the online supplementary material.

Supplementary Information The online version contains supplementary material available at https://doi.org/10.1007/s10463-021-00798-z.

Acknowledgements The authors sincerely thank the Editor, Associate Editor, and two anonymous reviewers for their insightful comments and suggestions. Sanying Feng's research was supported by the National Statistical Science Research Project of China (No. 2019LY18), the Natural Science Foundation of Henan Province, China (No. 212300410412), the Foundation of Henan Educational Committee, China (No. 21A910004), and the Excellent Youth Foundation of Zhengzhou University, China (No. 32210452). Tiejun Tong's research was supported by the General Research Fund (No. HKBU12303918), the National Natural Science Foundation of China (No. 1207010822), and the Initiation Grant for Faculty Niche Research Areas (No. RC-IG-FNRA/17-18/13) of Hong Kong Baptist University.

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