Two Sample Inference in Functional Linear Models

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Two sample inference in functional linear models

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Abstract

We propose a method of comparing two functional linear models in which explanatory variables are functions (curves) and responses can be either scalars or functions. In such models, the role of parameter vectors (or matrices) is played by integral operators acting on a function space. We test the null hypothesis that these operators are the same in two independent samples. The complexity of the test statistics increases as we move from scalar to functional responses and relax assumptions on the covariance structure of the regressors. They all, however, have an asymptotic chi-squared distribution with the number of degrees of freedom which depends on a specific setting. The test statistics are readily computable using the R package fda, and have good finite sample properties. The test is applied to egg-laying curves of Mediterranean flies and to data from terrestrial magnetic observatories.

Keywords: Functional linear model, significance test

Abbreviated Title: Two sample inference

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1 Introduction

In the last decade, methods of functional data analysis have been found useful in a number of applied fields, including environmental science, finance, genetics, biology, public health, geophysics, image and signal processing, to name just a few, see, for example, Besse et al. (2000), Damon and Guillas (2002), Laukaitis and Račkauskas (2002), Ando et al. (2004), Fernández de Castro et al. (2005), Morris et al. (2006), Müller et al. (2006), Hlubinka and Prchal (2007), Nerini and Ghattas (2007), Febrero et al. (2007), Glendinning and Fleet (2007). The functional linear model, in its various forms, stands out as a particularly useful tool in many such analyses, and has consequently been thoroughly studied, see Cuevas et al. (2002), Malfait and Ramsay (2003), Cardot et al. (2003a), Chiou et al. (2004), Müller and Stadtmüller (2005), Yao et al. (2005), Cai and Hall (2006), Chiou and Müller (2007), Li and Hsing (2007), among others.

This paper is concerned with the following problem of functional data analysis. We observe two samples: sample 1: \((X_i, Y_i), 1 \leq i \leq N\), and sample 2: \((X^*_j, Y^*_j), 1 \leq j \leq M\). The explanatory variables \(X_i\) and \(X^*_j\) are functions over a compact subset of a Euclidean space, whereas the responses \(Y_i\) and \(Y^*_j\) can be either functions or scalars (the \(Y_i\) and \(Y^*_j\) are either both functions, or both scalars). We model the dependence of the \(Y_i\) (\(Y^*_j\)) on the \(X_i\) (\(X^*_j\)) by the functional regression models

\[ Y_i = \Psi X_i + \varepsilon_i, \quad Y^*_j = \Psi^* X^*_j + \varepsilon^*_j, \]

where \(\Psi\) and \(\Psi^*\) are linear operators whose domain is a function space, and which take values either in the same function space or in the real line. We wish to test if the operators \(\Psi\) and \(\Psi^*\) are equal.

The above testing problem is directly motivated by our work with terrestrial magnetometer data. Specifically, we have been interested in finding out if interactions of currents of charged particles flowing in the ionosphere and the magnetosphere in various states of the near Earth space and at various times are statistically significantly different. The problem is truly functional, as the \(shapes\) of the daily magnetogram curves have been used for almost one hundred years to describe various electrodynamic phenomena.
in near Earth space. Digital observatories measure the components of the magnetic field every 5 seconds. In this application, which is described in some detail in Section 4, the model is fully functional (both responses and regressors are curves). The case of scalar responses is also included because of its importance in many applications, see e.g. Cardot et al. (2003a), Müller and Stadtmüller (2005) and Ferraty and Vieu (2006). Like classical two sample procedures in various forms, our methodology is likely to be applicable to a wide range of problems, where estimating two significantly different functional linear regressions on subsamples of a larger sample may reveal additional features.

Even in the traditional multivariate setting, except for comparing mean responses, the problem of comparing the regression coefficients for two models based on two different samples is not trivial, and we could not find a ready reference for it. In the functional setting, it becomes even more complex, as the dimension of the data must be reduced, and the resulting approximation error must be controlled. We achieve it by expanding the data with respect to the functional principal components (FPC’s), see e.g. Chapter 8 of Ramsay and Silverman (2005). In addition to being random, the FPC’s can be very different for the two samples, a problem not encountered in existing literature on functional linear models. Generally, due to different FPC’s structures, working with two functional samples is difficult, but important theoretical advances, mostly in the context of canonical correlation analysis, have recently been made, see He et al. (2003), Cupidon et al. (2007) and Eubank and Hsing (2007). Our paper aims at developing a practical testing procedure, but some complexity in the form of the test statistics and the derivation of their asymptotic properties is unavoidable. Our theory is developed using population, rather than estimated FPC’s. Estimation of the FPC’s introduces additional terms, and would make the arguments even longer. The additional terms are generally asymptotically or practically negligible, and the simulations we present here show that this is indeed the case in our setting. Gabrys and Kokoszka (2007), Horváth et al. (2009), Berkes et al. (2009), among others, show how the estimation of the FPC’s is handled in simpler settings.

Other approaches to comparing two samples of curves, but not samples of dependent pairs of curves, have been explored in Laukaitis and Račkauskas (2005), Ferraty et al.
(2007), Benko et al. (2009).

After introducing the requisite concepts and notation in Section 2, we describe the testing procedures in Section 3. Applications to medfly and magnetometer data are presented in Section 4. The asymptotic results and their proofs are collected in Section 5.

2 Preliminaries

We state the assumptions and introduce the notation for the observations \((X_i, Y_i), i = 1, 2, \ldots, N\). The observations \((X_j^*, Y_j^*), j = 1, 2, \ldots, M\), are assumed to satisfy the same assumptions with the superscript * added to the corresponding quantities. The two samples are assumed to be independent.

Throughout this paper we take \(\langle \cdot, \cdot \rangle\) to be the \(L^2([0, 1])\) inner product.

Assumption 2.1 The observations \(\{X_n\}\) are iid mean zero random functions in \(L^2([0, 1])\) satisfying

\[
\mathbb{E}||X_n||^4 = \mathbb{E} \left[ \int X_n^2(t) dt \right]^2 < \infty.
\]

For the linear model with scalar responses, we formulate the following assumption.

Assumption 2.2 The scalar responses \(Y_i\) satisfy

\[
Y_i = \int_0^1 \psi(s)X_i(s)ds + \varepsilon_i,
\]

with iid mean zero errors \(\varepsilon_i\) satisfying \(\mathbb{E}\varepsilon_i^4 < \infty\), and \(\psi \in L^2([0, 1])\). The errors \(\varepsilon_i\) and the regressors \(X_i\) are independent.

In the case of functional responses, we define an analogous assumption.

Assumption 2.3 The functional responses \(Y_i \in L^2([0, 1])\) satisfy

\[
Y_i(t) = \int_0^1 \psi(t, s)X_i(s)ds + \varepsilon_i(t),
\]
with iid mean zero errors $\varepsilon_i$ satisfying

$$E||\varepsilon_n||^4 = E \left[ \int \varepsilon_n^2(t)dt \right]^2 < \infty,$$

and $\psi \in L^2([0,1] \times [0,1])$. The errors $\varepsilon_i$ and the regressors $X_i$ are independent.

Our objective is to test

$$H_0 : \|\psi - \psi^*\| = 0$$

against

$$H_A : \|\psi - \psi^*\| \neq 0,$$

where the norm is in $L^2([0,1])$ under Assumption 2.2 and in $L^2([0,1] \times [0,1])$ under Assumption 2.3.

The covariance operator of the $X_n$ is defined by

$$C(x) = E[\langle X_n, x \rangle X_n], \quad x \in L^2([0,1])$$

and its eigenfunctions $v_j$ and eigenvalues $\lambda_j$ by $C(v_j) = \lambda_j v_j$, $j \geq 1$. The eigenfunctions $v_j$ form an orthonormal basis of $L^2([0,1])$. We also call them the functional principal components (FPC’s).

The empirical covariance operator is defined by

$$\hat{C}(x) = \frac{1}{N} \sum_{n=1}^{N} \langle X_n, x \rangle X_n, \quad x \in L^2([0,1]),$$

and its eigenelements by $\hat{C}(\hat{v}_j) = \hat{\lambda}_j \hat{v}_j$, $j = 1, 2, \ldots, N$.

Since the operators $C$ and $\hat{C}$ are symmetric and nonnegative definite, the eigenvalues $\lambda_j$ and $\hat{\lambda}_j$ are nonnegative. Following the usual convention, we assume that $\lambda_1 > \lambda_2 > \ldots$ with the same ordering for the empirical eigenvalues. Recall that $\sum_{j=1}^{\infty} \lambda_j = E||X_n||^2$, see e.g. Section 1.5 of Bosq (2000), so $v_1$ explains the largest portion of the variance of $X$, with the subsequent FPC’s explaining decreasing contributions to the variance. Our asymptotic results require that the first $p$ eigenvalues are nonzero and distinct, so we impose an additional assumption.

**Assumption 2.4** The eigenvalues of the covariance operator $C$ satisfy

$$\lambda_1 > \lambda_2 > \ldots \lambda_p > \lambda_{p+1}.$$
3 Testing procedure

In this section we introduce the testing procedure. We start with the easier case of scalar responses, and then generalize to the more technical case of functional response variables, which uses the results developed in the scalar setting. A reader interested only in the practical application can start with Section 3.3, which provides an outline.

3.1 Scalar responses

Since $\psi \in L^2([0,1])$, we can expand it as $\psi(s) = \sum_{i=1}^{\infty} \mu_i v_i(s)$, where $\mu_i = \langle \psi, v_i \rangle$. Consequently, the response variables can be expressed as $Y_i = \sum_{k=1}^{\infty} \mu_k \langle X_i, v_k \rangle + \varepsilon_i$. We truncate the above expansion at $1 \leq p < \infty$, and combine the error made by the truncation with the $\varepsilon_i$. Defining $\varepsilon'_i = \varepsilon_i + \sum_{k=p+1}^{\infty} \mu_k \langle X_i, v_k \rangle$, the response is equivalently given by

$$Y_i = \sum_{k=1}^{p} \mu_k \langle X_i, v_k \rangle + \varepsilon'_i. \quad (3.1)$$

In terms of matrix and vector notation we have

$$Y = X \mu + \varepsilon',$$

where, for $1 \leq i \leq N$ and $1 \leq j \leq p$,

$$Y(i) = Y_i, \quad X(i,j) = \langle X_i, v_j \rangle, \quad \mu(j) = \mu_j, \quad \varepsilon'(i) = \varepsilon'_i.$$

The least squares estimator for $\mu$ is therefore

$$\hat{\mu} = (X^T X)^{-1} X^T Y. \quad (3.2)$$

By Theorem 5.1, $\hat{\mu}$ is a consistent estimator of $\mu$, and for the second sample, the analogously defined $\hat{\mu}^*$ a consistent estimator of $\mu^*$. Thus we can base a test statistic on the difference $\hat{\mu} - \hat{\mu}^*$. To motivate our construction, assume first that the covariance operators of the $X_i$ and $X_j^*$ are equal and the errors $\varepsilon_i$ and $\varepsilon_j^*$ have equal variances, i.e.

$$E(X_1(s)X_1(t)) = E(X_1^*(s)X_1^*(t)) = c(s,t) \quad (3.3)$$

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and

\[ \text{var}(\varepsilon_1) = \text{var}(\varepsilon^*_1). \]  

(3.4)

In order to develop meaningful asymptotic results we assume that the two samples sizes are of roughly the same order. i.e. we assume there exists a constant \(0 < \zeta < \infty\) such that

\[ \frac{N}{M} \to \zeta, \quad N \to \infty. \]  

(3.5)

Finally we introduce the random variable

\[ \Lambda_p = N(1 + \zeta)^{-1} \left( \hat{\mu} - \hat{\mu}^* \right)^T \Sigma_p^{-1} \left( \hat{\mu} - \hat{\mu}^* \right), \]  

(3.6)

where \(\Sigma_p\) is the common asymptotic covariance matrix of \(\hat{\mu}\) and \(\hat{\mu}^*\) defined by

\[ \Sigma_p(i,i) = \lambda_i^{-1} \sigma^2 + \lambda_i^{-2} \text{var} \left( \langle X_1, v_i \rangle \sum_{k=p+1}^{\infty} \mu_k \langle X_1, v_k \rangle \right), \quad i = 1, \ldots, p; \]  

(3.7)

\[ \Sigma_p(i,j) = \lambda_i^{-1} \lambda_j^{-1} \text{E} \left( \langle X_1, v_i \rangle \langle X_1, v_j \rangle \left( \sum_{k=p+1}^{\infty} \mu_k \langle X_1, v_k \rangle \right)^2 \right), \quad i \neq j. \]  

(3.8)

The estimation of \(\Sigma_p\) and \(\Lambda_p\) is discussed later on in this section.

The following theorem, which follows immediately from Theorem 5.2, shows that \(\Lambda_p\) has a chi-squared asymptotic distribution.

**Theorem 3.1** Suppose Assumptions 2.1, 2.2, 2.4, and conditions (3.3), (3.4), and (3.5) hold. Then, \(\Lambda_p \overset{d}{\to} \chi^2(p)\), as \(N \to \infty\), where \(\Lambda_p\) is defined by (3.6), and \(\chi^2(p)\) is a chi-squared random variable with \(p\) degrees of freedom.

We therefore propose the following test statistic when the covariances are equal

\[ \hat{\Lambda}_p = N(1 + N/M)^{-1} \left( \hat{\mu} - \hat{\mu}^* \right)^T \Sigma_p^{-1} \left( \hat{\mu} - \hat{\mu}^* \right). \]  

(3.9)

We use the empirical diagonal approximation to the matrix \(\Sigma_p\) given by

\[ \Sigma_p = \hat{\sigma}^2 \begin{bmatrix} \hat{\lambda}_1^{-1} & 0 & \cdots & 0 \\ 0 & \hat{\lambda}_2^{-1} & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & \cdots & \hat{\lambda}_p^{-1} \end{bmatrix}; \]  

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The estimation of \(\Sigma_p\) and \(\Lambda_p\) is discussed later on in this section.
where \( \hat{\sigma} \) are residual standard deviations from the estimated regression models. The matrix \( \hat{\Sigma}_p \) is defined analogously.

In many applications, the covariance kernels \( c(s, t) \) and \( c^*(s, t) \) are not necessarily equal. Since the two kernels can have potentially different eigenfunctions we instead consider an arbitrary basis \( \{ w_i \} \) of \( L^2([0, 1]) \). The kernels \( \psi \) and \( \psi^* \) are expanded as

\[
\psi(s) = \sum_{i=1}^{\infty} \mu_i w_i(s), \quad \psi^*(s) = \sum_{j=1}^{\infty} \mu_j^* w_j(s),
\]

and so

\[
Y_i = \sum_{k=1}^{\infty} \mu_k \langle X_i, w_k \rangle + \varepsilon_i, \quad Y_i^* = \sum_{k=1}^{\infty} \mu_k^* \langle X_i^*, w_k \rangle + \varepsilon_i^*.
\]

Truncating both sums at \( p \), the response variables can again be expressed as

\[
Y = X\mu + \varepsilon', \quad Y^* = X^*\mu^* + \varepsilon'^*,
\]

with all terms analogously defined with respect to our new basis. While this appears similar to our prior calculations, we are expanding with respect to an arbitrary basis which means that \( X \) and \( \varepsilon' \) are now potentially correlated. The least squares estimators take the same form

\[
\hat{\mu} = (X^T X)^{-1} X^T Y, \quad \hat{\mu}^* = (X^{*T} X^*)^{-1} X^{*T} Y^*.
\]

Thus we can once again compare \( \hat{\mu} \) and \( \hat{\mu}^* \) to test the null hypothesis. To analyze the asymptotic behavior of these estimates we consider the relation

\[
\hat{\mu} = \mu + (X^T X)^{-1} X^T \varepsilon'.
\]

The vector \( X^T \varepsilon' \) can be expressed as

\[
X^T \varepsilon' = A + B + Nm,
\]

where

\[
A = \begin{pmatrix}
\sum_{i=1}^{N} \varepsilon_i \langle X_i, w_1 \rangle \\
\vdots \\
\sum_{i=p}^{N} \varepsilon_i \langle X_i, w_p \rangle
\end{pmatrix}, \quad B = \begin{pmatrix}
\sum_{i=1}^{N} \sum_{k=p+1}^{\infty} \mu_k (\langle X_i, w_1 \rangle \langle X_i, w_k \rangle - E[\langle X_1, w_1 \rangle \langle X_1, w_k \rangle]) \\
\vdots \\
\sum_{i=1}^{N} \sum_{k=p+1}^{\infty} \mu_k (\langle X_i, w_p \rangle \langle X_i, w_k \rangle - E[\langle X_1, w_p \rangle \langle X_1, w_k \rangle])
\end{pmatrix},
\]

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have mean zero and are uncorrelated since the error terms are independent of the ex-
planatory functions. The term \( m \) represents the bias introduced by using an arbitrary
basis which is given by

\[
m = \left( \sum_{k=p+1}^{\infty} \mu_k E[\langle X_1, w_1 \rangle \langle X_1, w_k \rangle] \begin{pmatrix} \vdots \\ \sum_{k=p+1}^{\infty} \mu_k E[\langle X_1, w_p \rangle \langle X_1, w_k \rangle] \end{pmatrix} \right).
\]

This yields the form

\[
\hat{\mu} = \mu + (X^T X)^{-1} A + (X^T X)^{-1} B + N(X^T X)^{-1} m.
\]

Clearly \( A \) and \( B \) are sums of iid random vectors with means zero and finite covariance
matrices due to Assumptions 2.1 and 2.2. Thus by the multivariate central limit theorem
\( N^{-1/2}(A \quad B)^T \) is asymptotically normal. We have by the strong law of large numbers
that

\[
N^{-1} \sum_{i=1}^{N} \langle X_i, w_j \rangle \langle X_i, w_k \rangle \xrightarrow{a.s.} E \langle X_1, w_j \rangle \langle X_1, w_k \rangle,
\]

for \( j = 1, \ldots, p \) and \( k = 1, \ldots, p \), or in matrix notation

\[
N^{-1} X^T X \xrightarrow{a.s.} \Sigma_1,
\]

where the \((j, k)\) entry of \( \Sigma_1 \) is \( E \langle X_1, w_j \rangle \langle X_1, w_k \rangle \). Thus by Slutsky’s Lemma \( N^{-1/2}(\hat{\mu} - \mu - N(X^T X)^{-1} m) \) is asymptotically normal. Since \( A \) has zero mean, we have that the
\((i, j)\) entry of its covariance matrix is given by

\[
E \sum_{k=1}^{N} \varepsilon_k \langle w_i, X_k \rangle \varepsilon_k \langle w_j, X_k \rangle = N \sigma^2 E \langle w_i, X_1 \rangle \langle w_j, X_1 \rangle,
\]

and therefore

\[
\text{cov}(A) = N \sigma^2 \Sigma_1.
\]

Turning to \( B \), the \((i, j)\) entry of its covariance matrix is given by

\[
N \sum_{k=p+1}^{\infty} \sum_{r=p+1}^{\infty} \mu_k \mu_r E \{ (\langle X_1, w_i \rangle \langle X_1, w_k \rangle - E[\langle X_1, w_i \rangle \langle X_1, w_k \rangle])
\]

\[
+ (\langle X_1, w_r \rangle \langle X_1, w_k \rangle - E[\langle X_1, w_r \rangle \langle X_1, w_k \rangle]) \}.
\]
\[(\langle X_1, w_j \rangle \langle X_1, w_r \rangle - E[\langle X_1, w_j \rangle \langle X_1, w_r \rangle])\].

We will denote the covariance matrix of \(B\) as \(N \Sigma_2\). Combining everything, we have by Slutsky’s Lemma

\[N^{1/2}(\hat{\mu} - \mu - N(X^T X)^{-1}m) \xrightarrow{d} N(0, C),\]

where \(C = \sigma^2 \Sigma_1^{-1} + \Sigma_1^{-1} \Sigma_2 \Sigma_1^{-1}\).

An identical argument gives, for the second sample,

\[M^{1/2}(\hat{\mu}^* - \mu^* - M(X^* T X^*)^{-1}m^*) \xrightarrow{d} N(0, C^*),\]

with all terms analogously defined. Combined with assumption 3.5 we therefore conclude

\[N^{1/2}(\hat{\mu} - \hat{\mu}^* - (N(X^T X)^{-1}m - M(X^* T X^*)^{-1}m^*)) \xrightarrow{d} N(0, C + \zeta C^*).\]

Neglecting the biases \(m\) and \(m^*\), we thus arrive at the test statistic

\[\hat{\Lambda}_p = N(\hat{\mu} - \hat{\mu}^*)^T (\hat{\sigma}^2 \hat{\Sigma}_1 + (N/M)\hat{\sigma}^*\hat{\Sigma}_1^*)^{-1}(\hat{\mu} - \hat{\mu}^*).\] (3.11)

The error variances \(\sigma^2\) and \(\sigma^*^2\) are estimated as before, but without pooling the samples. The matrix \(\Sigma_1\) is now estimated with

\[\hat{\Sigma}_1 = N^{-1}X^TX,\]

with \(\hat{\Sigma}_1^*\) defined analogously.

The distributions of statistics (3.9) and (3.11) are approximated by the chi-squared distribution with \(p\) degrees of freedom. If \(p\) is large (in terms of the percentage of variance explained), then all neglected terms are close to 0.

### 3.2 Functional responses

Turning to the case when the response variables are functions, we now assume that \(\varepsilon_i\) is a random element of \(L^2([0, 1])\) and \(\psi(t, s) \in L^2([0, 1] \times [0, 1])\) is fixed, but known. Our goal is to use dimension reduction techniques to make this problem approachable. Our
method will consist of choosing a basis to project the $Y_i$ onto, and then using the results developed in the scalar case.

We first focus on the case of equal variances defined by assumptions (3.3) and, in place of (3.4),

$$E(\varepsilon_1(s)\varepsilon_1(t)) = E(\varepsilon_1^*(s)\varepsilon_1^*(t)).$$

(3.12)

Consider an arbitrary orthonormal basis $\{u_i\}_{i=1}^{\infty}$ for $L^2([0,1])$ (on which the $Y_i$ are to be projected), and analogous basis $\{u^*_j\}_{j=1}^{\infty}$. Though all our results hold for an arbitrary choice for $\{u_i\}_{i=1}^{\infty}$, we will use in our applications the eigenfunctions of the covariance operator for the $\{Y_i\}$, with the $\{u^*_i\}$ defined analogously. Because $\{u_i\}$ and $\{v_i\}$ are both bases for $L^2([0,1])$, it follows that we can construct a basis for $L^2([0,1] \times [0,1])$ using the bivariate functions $u_i(t)v_j(s)$ for $(t,s) \in [0,1] \times [0,1]$, $i = 1, \ldots, \infty$, and $j = 1, \ldots, \infty$. We therefore have the expansion $\psi(t,s) = \sum_{k=1}^{\infty} \sum_{l=1}^{\infty} \mu_{kl} u_l(t) v_k(s)$, but we will work with the approximation

$$\hat{\psi}(t,s) = \sum_{k=1}^{p} \sum_{l=1}^{r} \hat{\mu}_{kl} \hat{u}_l(t) \hat{v}_k(s),$$

where $1 \leq r < \infty$ and $1 \leq p < \infty$ are fixed.

Extending the notation introduced in the case of scalar responses, define the matrices

$$Y(i,j) = \langle Y_i, u_j \rangle, \quad i = 1, \ldots, N, \quad j = 1, \ldots, r,$$
$$X(i,j) = \langle X_i, v_j \rangle, \quad i = 1, \ldots, N, \quad j = 1, \ldots, p,$$
$$\mu(i,j) = \int_0^1 \int_0^1 \psi(t,s) v_i(s) u_j(t) ds dt, \quad i = 1, \ldots, p, \quad j = 1, \ldots, r.$$

As in the model, we combine any errors made by our approximations with the error of the model, so we also introduce the matrix

$$\varepsilon'(i,j) = \langle \varepsilon_i, u_j \rangle + \sum_{k=p+1}^{\infty} X(i,k) \mu(k,j), \quad i = 1, \ldots, N, \quad j = 1, \ldots, r.$$

Projecting the relation $Y_i = \Psi X_i + \varepsilon_i$ onto the $u_j$, we obtain

$$\langle Y_i, u_j \rangle = \langle \Psi X_i, u_j \rangle + \langle \varepsilon_i, u_j \rangle = \sum_{k=1}^{\infty} \langle X_i, v_k \rangle \langle \Psi v_k, u_j \rangle + \langle \varepsilon_i, u_j \rangle$$
which implies
\begin{equation}
Y = X\mu + \varepsilon'.
\end{equation}

The corresponding least squares estimator \( \hat{\mu} = (X^TX)^{-1}X^TY \) consistently estimates the matrix \( \mu \). This follows immediately by applying Theorem 5.1 to each column of \( \hat{\mu} \). Asymptotic normality follows from Theorem 5.4.

Since \( \mu \) is now a matrix, the task of constructing a quadratic form leading to a test statistic is somewhat painful notationally. We start by writing \( \mu \) as a column vector of length \( pr \):
\begin{equation}
\mu_v^T = \text{vec}(\mu) = (\mu(1,1), \mu(2,1), \ldots, \mu(p,1), \mu(1,2), \ldots, \mu(p-1,r), \mu(p,r)).
\end{equation}

In words, \( \mu_v \) is constructed by placing the columns of \( \mu \) on top of one another. The covariance matrix for the error terms is given by
\[ \Sigma_{\varepsilon}(i,j) = \text{cov}[\langle \varepsilon_1, u_i \rangle, \langle \varepsilon_1, u_j \rangle] = E[\langle \varepsilon_1, u_i \rangle \langle \varepsilon_1, u_j \rangle], \quad 1 \leq i, j \leq r, \]
and the diagonal matrix containing the largest \( p \) eigenvalues of \( C \) is
\[ \Gamma(i,j) = \lambda_i \delta_{ij}, \quad \text{for } 1 \leq i, j \leq p, \]
where \( \delta_{ij} \) is Kronecker's delta.

With this notation in place, we consider the random variable
\[ \Lambda_{pr} = N(1 + \zeta)^{-1}(\hat{\mu}_v - \mu_v^*)^T \left( \Sigma_{\varepsilon} \otimes \Gamma^{-1} + E \left[ \Delta_1 \otimes (\Gamma^{-1} \Delta_2 \Gamma^{-1}) \right] \right)^{-1} (\hat{\mu}_v - \mu_v^*), \]
where \( \Delta_1, \Delta_2 \) are defined in (5.4) and (5.5), respectively.

Assuming equal covariances, Theorem 5.4 implies that under \( H_0, \Lambda_{pr} \xrightarrow{d} \chi^2(pr) \). An extension of the argument used in the proof of Theorem 3.1 yields that \( \Lambda_{pr} \xrightarrow{p} \infty \), under \( H_A \), as long as \( p \) and \( r \) are so large that \( \mu \neq \mu^* \). That such a pair \( (p,r) \) exists, follows immediately from the fact that \( \{v_i u_j\} \) form a basis in \( L_2([0,1] \times [0,1]) \).

A computable approximation to \( \Lambda_{pr} \) is
\begin{equation}
\hat{\Lambda}_{pr} = N(1 + N/M)^{-1}(\hat{\mu}_v - \mu_v^*)^T \left( \hat{\Sigma}_{\varepsilon} \otimes \hat{\Gamma}^{-1} \right)^{-1} (\hat{\mu}_v - \mu_v^*),
\end{equation}
where $\Sigma_\varepsilon$ is the sample covariance matrix of the residuals in (3.13) and $\hat{\Gamma} = \text{diag}(\hat{\lambda}_1, \hat{\lambda}_2, \ldots, \hat{\lambda}_p)$, with the $\hat{\lambda}_i$ being the eigenvalues of the empirical covariance operator of the pooled $X_i$ and $X_j^*$.

We finally turn to the most complex case of different covariances for the explanatory functions. We now expand both the explanatory and response functions with respect to two arbitrary, potentially different, bases in $L^2[0, 1]$, $\{u_i\}$ and $\{w_j\}$, respectively:

$$
\psi(t, s) = \sum_{i=1}^{\infty} \sum_{j=1}^{\infty} \mu_{ji} u_i(t) w_j(s), \quad \psi^*(t, s) = \sum_{i=1}^{\infty} \sum_{j=1}^{\infty} \mu^*_{ji} u_i(t) w_j(s).
$$

This leads to the relations $Y = X\mu + \varepsilon'$, $Y^* = X^*\mu^* + \varepsilon'^*$, with all terms analogously defined as in the equal variance case, but using the bases $\{u_i\}$ and $\{w_j\}$. Thus the least squares estimates are again $\hat{\mu} = (X^TX)^{-1}X^TY$ and $\hat{\mu}^* = (X^{*T}X^*)^{-1}X^{*T}Y^*$. $(\mu, \mu^*, \hat{\mu}, \hat{\mu}^*)$ are now $p \times r$ matrices) Extending the argument developed in Section 3.1, we arrive at the test statistic

$$
\hat{\Lambda}_{pr} = N\text{vec}(\hat{\mu} - \hat{\mu}^*)^T (\hat{\Sigma}_\varepsilon \otimes \hat{\Sigma}_1^{-1} + (N/M)\hat{\Sigma}_\varepsilon^* \otimes \hat{\Sigma}_1^{*-1})^{-1}\text{vec}(\hat{\mu} - \hat{\mu}^*),
$$

where the residual covariance matrices $\Sigma_\varepsilon$ and $\Sigma_\varepsilon^*$ are computed for each sample separately. The estimate $\hat{\Sigma}_1^{-1}$ is given by $N^{-1}X^TX$, and $\hat{\Sigma}_1^{*-1}$ is equivalently defined.

The distribution of statistics (3.15) and (3.16) is approximated by the chi-square distribution with $pr$ degrees of freedom. The values of $p$ and $r$ depend on the selected bases, examples of suitable bases are given in Section 4. Selection of $p$ and $r$ is discussed in Section 3.3.

If $p$ and/or $r$ are large, the normalized $\chi^2$ distribution can be approximated by a normal distribution, as in Cardot et al. (2003a) who studied a single scalar response model and tested $\psi = 0$. In our case, due to the complexity of the problem, the rigorous derivation of the normal convergence with $p = p_n$ depending on a sample size would be far more tedious, so it is not pursued. To perform a test, a finite $p$ (and $r$) must be chosen no matter what approximation is used, and as illustrated in Section 4 large $p$ (and $r$) do not necessarily lead to meaningful results.
3.3 Summary of the testing procedure

In order to apply the tests, we must first verify if a linear functional model approximates the dependence structure of the data reasonably well. This can be done using the techniques of Chiou and Müller (2007). The assumptions of independence and identical distribution can be verified using the test of Gabrys and Kokoszka (2007). The regressors and the responses must be centered, so that their sample means are zero.

Next, the values of $p$ and $r$ must be chosen. In applications in which the FPC’s have a clear interpretation, these values can be chosen so that the action of the operators on specific subspaces spanned by the FPC’s of interest is compared. In the absence of such an interpretation, several data driven approaches are available. When the covariances are approximately equal, typically $p$ is chosen so large that $\sum_{k=1}^{p} \hat{\lambda}_k$ exceeds a required percentage of the variance of the $X_i$ (defined as $(N+M)^{-1}(\sum_{i=1}^{N} \int X_i^2(t)dt + \sum_{j=1}^{M} \int X_i^*2(t)dt)$ for the centered functions). We choose $r$ analogously for the response functions. When the covariances cannot be assumed equal then we propose, as one possibility, a pooling technique to choose $p$ and $r$. Pooling the explanatory functions we have

$$(N + M)^{-1} \left( \sum_{i=1}^{N} X_i(s)X_i(t) + \sum_{j=1}^{M} X_j^*(s)X_j^*(t) \right) \xrightarrow{a.s.} (1 + 1/\zeta)^{-1}c(s,t) + (1 + \zeta)c^*(s,t).$$

We propose taking the $w_i$ to be the eigenfunctions of $(1 + 1/\zeta)^{-1}c(s,t) + (1 + \zeta)c^*(s,t)$ which is itself a covariance kernel. The $u_i$ can be defined in an analogous manner using the response functions. Such a choice will allow smaller values of $p$ (and $r$) to be taken so that any bias from neglected terms is minimal, but we can still expect reasonable power. The values $p$ and $r$ can be chosen as before, but now with respect to the pooled variance.

All these steps can be implemented in the R package `fda`, and ready–made functions for the percentage of variance explained by FPC’s are available. Other methods of choosing $p$ (or $r$) are implemented in the MATLAB PACE package developed at the University of California at Davis.

It is often useful to compute the test for a wide range of values of $p$ (and $r$) and check if a uniform pattern emerges. This approach is illustrated in Section 4.
Finally, we compute the test statistic $\hat{\Lambda}$, and reject $H_0$ if it exceeds the $\chi^2$ density with DF degrees of freedom according to the following table:

<table>
<thead>
<tr>
<th>Response</th>
<th>Covariances</th>
<th>$\hat{\Lambda}$</th>
<th>DF</th>
</tr>
</thead>
<tbody>
<tr>
<td>Scalar</td>
<td>Equal</td>
<td>(3.9)</td>
<td>$p$</td>
</tr>
<tr>
<td>Scalar</td>
<td>Different</td>
<td>(3.11)</td>
<td>$p$</td>
</tr>
<tr>
<td>Functional</td>
<td>Equal</td>
<td>(3.15)</td>
<td>$pr$</td>
</tr>
<tr>
<td>Functional</td>
<td>Different</td>
<td>(3.16)</td>
<td>$pr$</td>
</tr>
</tbody>
</table>

The term “equal covariances” refers to assumptions (3.3), (3.4) in the scalar case, and (3.3), (3.12) in the functional case.

4 Application to medfly and magnetometer data

In this section we illustrate the application of the test on two examples. The first example is motivated by the work presented in Carey et al. (2002), Chiou et al. (2004), Müller and Stadtmüller (2005), Chiou and Müller (2007), among others, and studies egg-laying curves of Mediterranean fruit flies (medflies). The second example is an application to the measurements of the magnetic field generated by near Earth space currents. Techniques of FDA are very promising in space physics, see Kokoszka et al. (2008) and Maslova et al. (2009).

Before turning to these examples, we note that simulations performed on generic models show that the procedures have empirical sizes are close to nominal, and the power increases with the “size” of the difference. The power is smaller if the explanatory functions do not have a common distribution, and/or are non-normal. Some tables and the R code are available at http://www.stat.uchicago.edu/~mreimherr/research.html.

**Egg-laying curves of Mediterranean fruit flies.** Müller and Stadtmüller (2005), Section 6, consider 534 egg-laying curves of medflies who lived at least 30 days. Each function is defined over an interval $[0, 30]$, and its value on day $t \leq 30$ is the count of eggs laid by fly $i$ on that day. The 534 flies can be classified into long–lived, i.e. those
who lived longer than 44 days, and short–lived, i.e. those who died before the end of
the 44th day after birth. In the sample, there are 256 short–lived, and 278 long–lived
flies. This classification naturally defines two samples: Sample 1: the egg-laying curves
$X_i(t), \ 0 < t \leq 30, \ i = 1, 2, \ldots, 256$ of the short–lived flies, and the corresponding total
number of eggs $Y_i$. Sample 2: the egg-laying curves $X^*_j(t), \ 0 < t \leq 30, \ j = 1, 2, \ldots, 278$
of the long–lived flies, and the corresponding total number of eggs $Y^*_j$.

The average of the $Y^*_j$ is obviously larger than that of the $Y_i$, but a question of
interest is whether after adjusting for the means, the structure of the dependence of the
$Y^*_j$ on the curves $X^*_j(t)$ is different from the dependence of the $Y_i$ on the curves $X_i(t)$. Additionally, we would like to know in which functional component the difference, if any,
becomes significant. As the graphs and the analysis in Müller and Stadtmüller (2005)
show, even distinguishing the two functional samples $X_i(t), \ 0 < t \leq 30, \ i = 1, 2, \ldots, 256$
and $X^*_j(t), \ 0 < t \leq 30, \ j = 1, 2, \ldots, 278$ is very difficult.

We thus consider two linear models:

$$ Y_i - \bar{Y} = \int \psi(t)(X_i(t) - \bar{X}(t))dt + \varepsilon_i, \ i = 1, 2, \ldots, 256, $$

$$ Y^*_j - \bar{Y}^* = \int \psi^*(t)X^*_j(t) - \bar{X}^*(t))dt + \varepsilon^*_j, \ j = 1, 2, \ldots, 278. $$

and test (without assuming equal variances) $H_0: \psi = \psi^*$.

The above linear models describe the dependence structure of the data remarkably
well. We applied the graphical test of Chiou and Müller (2007) in which the responses are
graphed against the scores of the initial functional principal components. Figure 1 shows
such graphs for the second principal component, the graphs for the other components also
show nice elliptic shapes.

Table 1 shows the P-values for the five initial FPC’s. The P-values for the remaining
FPC’s do not exceed half a percent. We see that there is a significant difference between
the curves $\psi$ and $\psi^*$, and that it occurs in the dependence on the second principal com-
ponents of the egg-laying curves. Using only the first principal components ($p = p^* = 1$)
is not enough to see the difference.

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Data from terrestrial magnetic observatories. We now apply our methodology to magnetometer data. A comprehensive case study is not our goal, but we would rather like to illustrate the steps outlined in Section 3.3 in a new, practically relevant setting.

A sample of 40 functional regressors and corresponding responses is shown in Figure 2. Each curve in Figure 2 shows one minute averages in a UT (Universal Time) day of the component of the magnetic field lying in the Earth’s tangent plane and pointing toward the magnetic North. We thus have 1440 data points per curve, so a traditional multivariate analysis is not feasible. However, splitting the magnetogram data into days and treating the daily curves as functional observations is natural because of the daily rotation of the Earth. Moreover, for space physics researchers, it is the shape of these curves that conveys information, not the minute by minute values. These curves record magnetic fields generated by the Magnetosphere–Ionosphere (M–I) current system. The M–I currents are located in a region of space which is not sufficiently dense to allow operation of aircraft, but too dense for satellites, so in situ measurements are difficult to
Table 1: Statistic (3.11) and the P-values for several values of $p = p^\ast$.

<table>
<thead>
<tr>
<th>$p$</th>
<th>$\Lambda_{pp^\ast}$</th>
<th>P-Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>1.3323</td>
<td>0.2484</td>
</tr>
<tr>
<td>2</td>
<td>11.3411</td>
<td>0.0034</td>
</tr>
<tr>
<td>3</td>
<td>10.6097</td>
<td>0.0140</td>
</tr>
<tr>
<td>4</td>
<td>23.8950</td>
<td>0.0001</td>
</tr>
<tr>
<td>5</td>
<td>33.1144</td>
<td>0.0000</td>
</tr>
</tbody>
</table>

obtain. For this reason, and because of its inherent complexity, the M–I current system is only partially understood. An interested reader is referred to Kivelson and Russell (1997), especially Chapters 9, 10, 13 and 14. Here we utilize a simple and easy to explain setting, intended to illustrate our technique. The curves $X_i$ reflect ionospheric magnetic activity in the polar region known as substorms, which is spectacularly manifested as northern lights (aurora borealis). The curves $Y_i$ reflect magnetospheric activity in the magnetic equatorial region in the same UT day. We consider three samples: A, B, C. Each of them consists of about 40 pairs of curves. All measurements were recorded in 2001, the $X_i$ at College (CMO), Alaska; the $Y_i$ at Honolulu (HON), Hawaii. Sample A contains substorms which took place in January through March, B in April–June, C in July–September. Broader physical issues related to the dependence of the $Y_i$ on the $X_i$ in Kamide et al. (1998). Using a goodness–of–fit test of Chiou and Müller (2007), Kokoszka et al. (2008) verified that the fully functional linear model is a reasonable approximation to the dependence of the $Y_i$ on the $X_i$. The identical distribution and lack of correlations of the observations were established in Gabrys and Kokoszka (2007) and Kokoszka et al. (2008). The data can be assumed to be approximately independent because the M-I system resets itself after each rotation of the Earth, and the effect of larger disturbances of solar origin decay within about two days.

Intuitively, we would expect rejections of the null for all three pairs: A–B, B–C, and
Figure 2: Observations for sample A: left panel CMO (X), right panel HON (Y).

A–C, as the position of the axis of the Earth relative to the Sun shifts with each season, and substorms are influenced by the solar wind. This is indeed the case for tests in cases B–C and A–C, for which the P-values are very small: for B–C the largest P-value is 0.034, and for A–C 0.007 (for $p \leq 15$, $r \leq 10$). The results for testing samples A and B presented in Table 2 indicate the acceptance of $H_0$. In retrospect, this conclusion is supported by the observation, well-known in the space-physics community, that M–I disturbances tend to be weaker in summer months. Our test thus shows that it is reasonable to assume that the effect of substorms on low-latitude currents is approximately the same in first and second quarter of 2001, but changes in the third quarter (due to weaker substorms).
Table 2: P–values for testing the equality of regression operators in samples A and B.

<table>
<thead>
<tr>
<th>p/r</th>
<th>1</th>
<th>2</th>
<th>3</th>
<th>4</th>
<th>5</th>
<th>6</th>
<th>7</th>
<th>8</th>
<th>9</th>
<th>10</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>0.344</td>
<td>0.608</td>
<td>0.231</td>
<td>0.349</td>
<td>0.380</td>
<td>0.372</td>
<td>0.391</td>
<td>0.351</td>
<td>0.257</td>
<td></td>
</tr>
<tr>
<td>2</td>
<td>0.147</td>
<td>0.259</td>
<td>0.274</td>
<td>0.416</td>
<td>0.565</td>
<td>0.422</td>
<td>0.373</td>
<td>0.345</td>
<td>0.339</td>
<td>0.310</td>
</tr>
<tr>
<td>3</td>
<td>0.204</td>
<td>0.378</td>
<td>0.399</td>
<td>0.621</td>
<td>0.592</td>
<td>0.582</td>
<td>0.621</td>
<td>0.654</td>
<td>0.478</td>
<td></td>
</tr>
<tr>
<td>4</td>
<td>0.120</td>
<td>0.305</td>
<td>0.299</td>
<td>0.567</td>
<td>0.716</td>
<td>0.619</td>
<td>0.654</td>
<td>0.307</td>
<td>0.315</td>
<td>0.158</td>
</tr>
<tr>
<td>5</td>
<td>0.440</td>
<td>0.668</td>
<td>0.555</td>
<td>0.741</td>
<td>0.861</td>
<td>0.730</td>
<td>0.792</td>
<td>0.515</td>
<td>0.453</td>
<td>0.223</td>
</tr>
<tr>
<td>6</td>
<td>0.582</td>
<td>0.891</td>
<td>0.798</td>
<td>0.793</td>
<td>0.883</td>
<td>0.554</td>
<td>0.567</td>
<td>0.605</td>
<td>0.218</td>
<td>0.106</td>
</tr>
<tr>
<td>7</td>
<td>0.689</td>
<td>0.962</td>
<td>0.950</td>
<td>0.911</td>
<td>0.954</td>
<td>0.749</td>
<td>0.792</td>
<td>0.783</td>
<td>0.566</td>
<td>0.427</td>
</tr>
<tr>
<td>8</td>
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<td>0.968</td>
<td>0.972</td>
<td>0.952</td>
<td>0.958</td>
<td>0.815</td>
<td>0.755</td>
<td>0.582</td>
<td>0.432</td>
<td>0.257</td>
</tr>
<tr>
<td>9</td>
<td>0.981</td>
<td>0.804</td>
<td>0.962</td>
<td>0.980</td>
<td>0.972</td>
<td>0.821</td>
<td>0.837</td>
<td>0.753</td>
<td>0.722</td>
<td>0.456</td>
</tr>
<tr>
<td>10</td>
<td>0.727</td>
<td>0.585</td>
<td>0.903</td>
<td>0.973</td>
<td>0.986</td>
<td>0.972</td>
<td>0.973</td>
<td>0.941</td>
<td>0.935</td>
<td>0.626</td>
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<td>0.998</td>
<td>0.998</td>
<td>0.994</td>
<td>0.995</td>
<td>0.990</td>
</tr>
<tr>
<td>12</td>
<td>0.856</td>
<td>0.860</td>
<td>0.989</td>
<td>0.997</td>
<td>0.959</td>
<td>0.962</td>
<td>0.940</td>
<td>0.930</td>
<td>0.845</td>
<td>0.889</td>
</tr>
<tr>
<td>13</td>
<td>0.667</td>
<td>0.856</td>
<td>0.982</td>
<td>0.988</td>
<td>0.939</td>
<td>0.950</td>
<td>0.889</td>
<td>0.845</td>
<td>0.784</td>
<td>0.844</td>
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<tr>
<td>14</td>
<td>0.395</td>
<td>0.457</td>
<td>0.798</td>
<td>0.418</td>
<td>0.314</td>
<td>0.445</td>
<td>0.240</td>
<td>0.240</td>
<td>0.201</td>
<td>0.282</td>
</tr>
<tr>
<td>15</td>
<td>0.398</td>
<td>0.481</td>
<td>0.847</td>
<td>0.414</td>
<td>0.321</td>
<td>0.456</td>
<td>0.276</td>
<td>0.255</td>
<td>0.170</td>
<td>0.113</td>
</tr>
</tbody>
</table>
5 Asymptotic theory

In this section we collect the statements and proofs of the theorems referred to in Section 3, and continue to use the notation introduced in that section. These results do not follow from the existing multivariate theory because the regression errors are not independent and include projections on the “left over” FPC’s $v_{p+1}, v_{p+2}, \ldots, u_{r+1}, u_{r+2}, \ldots, \text{etc.}$ Theorems 5.2 and 5.4 are of particular interest, as they state the exact asymptotic distribution of the LSE’s in a multivariate regression obtained by projecting a functional regression.

Since the following simple lemma is used repeatedly in the proofs, it is stated first for ease of reference. It essentially follows from the orthonormality of the eigenfunctions of the covariance operator.

**Lemma 5.1** Suppose $X$ is a random element of $L^2([0,1])$ satisfying

$$EX(t) = 0 \quad \text{and} \quad E \int_0^1 |X(t)|^2 dt < \infty.$$ 

Let $\lambda_i, v_i, \ i \geq 1,$ be, respectively, the eigenvalues and the eigenfunctions of its covariance operator. Then

$$E[\langle v_i, X \rangle \langle v_j, X \rangle] = \lambda_i \delta_{ij},$$

where $\delta_{ij}$ is Kronecker’s delta.

**Theorem 5.1** Suppose Assumptions 2.1, 2.2 and 2.4 hold. Then,

$$\hat{\mu} = (X^TX)^{-1} X^T Y \xrightarrow{a.s.} \mu, \quad \text{as } N \to \infty,$$

where $\mu^T = (\mu_1, \ldots, \mu_p)$ and $\xrightarrow{a.s.}$ refers to almost sure convergence.

**Proof:** To analyze the behavior of $\hat{\mu}$, let us start by considering

$$(X^TX)(i,j) = \sum_{k=1}^N \langle v_i, X_k \rangle \langle v_j, X_k \rangle.$$ 

Since the $X_i$ are iid, by the strong law of large numbers

$$\frac{1}{N}(X^TX)(i,j) \xrightarrow{a.s.} E(\langle v_i, X_1 \rangle \langle v_j, X_1 \rangle), \quad \text{as } N \to \infty.$$
From Lemma 5.1 we have that $E(\langle v_i, X_1 \rangle \langle v_j, X_1 \rangle) = \lambda_i \delta_{ij}$. Therefore $N^{-1}(X^T X)$ converges almost surely to a $p \times p$ diagonal matrix whose diagonal entries are the eigenvalues of $C$.

Turning to $X^T Y$, observe that

$$X^T Y(i) = \sum_{j=1}^{N} \langle v_i, X_j \rangle Y_j$$

and, using (3.1),

$$\sum_{j=1}^{N} \langle v_i, X_j \rangle Y_j = \sum_{j=1}^{N} \langle v_i, X_j \rangle \sum_{k=1}^{p} \mu_k \langle v_k, X_j \rangle + \sum_{j=1}^{N} \epsilon_j \langle v_i, X_j \rangle.$$

Applying again the strong law of large numbers and Lemma 5.1 again, we obtain, as $N \to \infty$,

$$N^{-1} \sum_{j=1}^{N} \langle v_i, X_j \rangle \sum_{k=1}^{p} \mu_k \langle v_k, X_j \rangle \overset{a.s.}{\to} E \sum_{k=1}^{p} \mu_k \langle v_i, X_1 \rangle \langle v_k, X_1 \rangle = \mu_i \lambda_i \delta_{ij}.$$

Lastly, we will show that, as $N \to \infty$,

$$N^{-1} \sum_{j=1}^{N} \epsilon_j \langle v_i, X_j \rangle \overset{a.s.}{\to} 0.$$

Recalling the definition of $\epsilon'_i$, (3.1), we have

$$N^{-1} \sum_{j=1}^{N} \epsilon'_j \langle v_i, X_j \rangle = N^{-1} \sum_{j=1}^{N} \epsilon_j \langle v_i, X_j \rangle + N^{-1} \sum_{j=1}^{N} \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_j \rangle \langle v_i, X_j \rangle.$$

Since $\{\epsilon_i\}$ and $\{X_i\}$ are independent, by the strong law of large numbers and Assumption 2.2

$$N^{-1} \sum_{j=1}^{N} \epsilon_j \langle v_i, X_j \rangle \overset{a.s.}{\to} 0.$$

Similarly, using Lemma 5.1 and noting that $i \leq p$, we get

$$N^{-1} \sum_{j=1}^{N} \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_j \rangle \langle v_i, X_j \rangle \overset{a.s.}{\to} E \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \langle v_i, X_1 \rangle = 0.$$

**Theorem 5.2** Suppose Assumptions 2.1, 2.2 and 2.4 hold. Then, as $N \to \infty$,

$$\sqrt{N}(\hat{\mu} - \mu) \overset{d}{\to} N(0, \Sigma_p),$$

where $N(0, \Sigma_p)$ is a multivariate normal random vector with mean 0 and covariance matrix $\Sigma_p$ defined by (3.7) and (3.8).
\textbf{Proof:} By the definition of $\hat{\mu} (3.2)$,
\[
\sqrt{N}(\hat{\mu} - \mu) = \sqrt{N} \left( \left( X^T X \right)^{-1} X^T Y - \mu \right).
\]
Defining $\epsilon'^T = (\epsilon'_1, \ldots, \epsilon'_N)$, the above reduces to
\[
\sqrt{N}(\hat{\mu} - \mu) = \left( N^{-1} X^T X \right)^{-1} N^{-1/2} X^T \epsilon'.
\]
By Lemma 5.1, $N^{-1} X^T X$ converges almost surely to a diagonal matrix whose diagonal elements are the first $p$ eigenvalues of the covariance operator of the $\{X_i\}$. Therefore we need only focus on the behavior of $N^{-1/2} X^T \epsilon'$ and use Slutsky’s Theorem to obtain the claimed limiting distribution. Considering the $i^{th}$ coordinate of $N^{-1/2} X^T \epsilon'$ we have
\[
(N^{-1/2} X^T \epsilon')(i) = N^{-1/2} \sum_{j=1}^{N} \langle v_i, X_j \rangle \epsilon'_j, \quad 1 \leq i \leq p.
\]
By Assumption 2.2 the above is a summation of iid random variables. Since each coordinate of $N^{-1/2} X^T \epsilon'$ is given by such a sum, Assumption 2.2 implies that $X^T \epsilon'$ can be expressed as a sum of iid random vectors. We can apply the multivariate central limit theorem to obtain the claimed multivariate normal limiting distribution if we can show that each entry of the covariance matrix is finite. Therefore we spend the rest of the proof deriving the form for $\Sigma_p$ and showing that its entries are finite. Using the definition of $\epsilon'_i$, we obtain
\[
N^{-1/2} \sum_{j=1}^{N} \langle v_i, X_j \rangle \epsilon'_j = N^{-1/2} \left( \sum_{j=1}^{N} \langle v_i, X_j \rangle \epsilon_j + \sum_{j=1}^{N} \langle v_i, X_j \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_j \rangle \right).
\]
Because the $\{X_j\}$ are independent, both sums (with respect to $j$) are sums of independent and identically distributed random variables. Furthermore, since $\{\epsilon_j\}$ are independent of all other terms, we also have that the two sums above are uncorrelated. Therefore it follows that
\[
\text{var} \left( N^{-1/2} \left( \sum_{j=1}^{N} \langle v_i, X_j \rangle \epsilon_j + \sum_{j=1}^{N} \langle v_i, X_j \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_j \rangle \right) \right)
\]
\[
= \text{var} \left( \langle v_i, X_1 \rangle \epsilon_1 \right) + \text{var} \left( \langle v_i, X_1 \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right).
\]
\[
(5.2)
\]
Considering the first term of (5.2), we have by the independence of $X_1$ and $\varepsilon_1$ and Lemma 5.1 that
\[
\text{var}(\langle v_i, X_1 \rangle \varepsilon_1) = \lambda_i \sigma^2 < \infty.
\]

Turning to the second term of (5.2), we have by Lemma 5.1
\[
\text{var} \left[ \left\langle v_i, X_1 \right\rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right] = \mathbb{E} \left[ \left\langle v_i, X_1 \right\rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^2.
\]

Applying the Cauchy-Schwarz inequality it follows that
\[
\mathbb{E} \left[ \left\langle v_i, X_1 \right\rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^2 \leq \left( \mathbb{E} \left[ \left\langle v_i, X_1 \right\rangle \right]^4 \mathbb{E} \left[ \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^4 \right)^{1/2}.
\]

As a consequence of Assumption 2.1, we obtain that
\[
\mathbb{E} \left[ \langle v_i, X_1 \rangle \right]^4 < \infty.
\]

Using the Cauchy-Schwarz inequality again we have
\[
\mathbb{E} \left[ \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^4 \leq \mathbb{E} \left[ \sum_{k=p+1}^{\infty} \mu_k^2 \sum_{s=p+1}^{\infty} \langle v_s, X_1 \rangle^2 \right]^2.
\]

Therefore we can infer that
\[
\mathbb{E} \left[ \langle v_i, X_1 \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^2 \leq \left( \sum_{k=p+1}^{\infty} \mu_k^2 \right)^{1/2} \left( \mathbb{E} \left[ \langle v_i, X_1 \rangle \right]^4 \mathbb{E} \left[ \sum_{s=p+1}^{\infty} \langle v_s, X_1 \rangle^2 \right]^4 \right)^{1/2}.
\]

Using Assumption 2.2 and Bessel’s Inequality we have that
\[
\sum_{k=p+1}^{\infty} \mu_k^2 \leq \| \psi \|^2 < \infty.
\]

Similarly, using Assumption 2.1 and Bessel’s Inequality we have that
\[
\mathbb{E} \left( \sum_{s=p+1}^{\infty} \langle v_s, X_1 \rangle^2 \right)^2 \leq \mathbb{E} \| X_1 \|^4 < \infty.
\]

Combining the above with Assumption 2.1 we conclude
\[
\mathbb{E} \left[ \langle v_i, X_1 \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right]^2 < \infty.
\]
and it follows that the diagonal elements of $\Sigma_p$ are given

$$\Sigma_p(i, i) = \lambda_i^{-1} \sigma^2 + \lambda_i^{-2} \text{var} \left( \langle v_i, X_1 \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right), \ i = 1, \ldots, p.$$ 

Next we examine the joint behavior of the coordinates. Combining (5.3) with the Cauchy-Schwarz inequality we have

$$\text{cov} \left[ (X^T \varepsilon')(i), (X^T \varepsilon')(j) \right] < \infty \quad i = 1, \ldots, p \quad \text{and} \quad j = 1, \ldots, p.$$ 

Therefore to finish the proof we need only derive the form for the off diagonal terms of $\Sigma_p$. Using (5.1), Assumption 2.2, and Lemma 5.1, it is easy to verify that for $i \neq j$

$$\text{cov} \left[ (X^T \varepsilon')(i), (X^T \varepsilon')(j) \right] = E((X^T \varepsilon')(i)(X^T \varepsilon')(j))$$

$$= E \left( \sum_{q=1}^{N} \langle v_i, X_q \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_q \rangle \sum_{s=1}^{N} \langle v_j, X_s \rangle \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_s \rangle \right)$$

$$= \sum_{q=1}^{N} E \left( \langle v_i, X_q \rangle \langle v_j, X_q \rangle \left( \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_q \rangle \right)^2 \right)$$

$$= NE \left( \langle v_i, X_1 \rangle \langle v_j, X_1 \rangle \left( \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right)^2 \right).$$ 

Therefore it follows that the off diagonal terms of $\Sigma_p$ are given by

$$\Sigma_p(i, j) = \lambda_i^{-1} \lambda_j^{-1} E \left( \langle v_i, X_1 \rangle \langle v_j, X_1 \rangle \left( \sum_{k=p+1}^{\infty} \mu_k \langle v_k, X_1 \rangle \right)^2 \right), \ i \neq j,$$

which concludes the proof.

**Theorem 5.3** If the assumptions of Theorem 3.1 are satisfied and $p$ is so large that $\mu \neq \mu^*$, then $\Lambda_p \overset{p}{\rightarrow} \infty$, as $N \rightarrow \infty$.

**Proof:** We start by expanding $\Lambda_p$ as

$$\Lambda_p = N(1 + \zeta)^{-1} (\hat{\mu} - \mu^*)^T \Sigma_p^{-1} (\hat{\mu} - \mu^*)$$

$$= N(1 + \zeta)^{-1} (\hat{\mu} - \mu - \hat{\mu}^* + \mu^*)^T \Sigma_p^{-1} (\hat{\mu} - \mu - \hat{\mu}^* + \mu^*)$$

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\[ N(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\mu - \mu^*) \]
\[ + 2N(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\hat{\mu} - \mu^* + \mu^*) . \]

Therefore we need only consider each term above. From Theorem 3.1 it follows that
\[ N(1 + \zeta)^{-1}(\hat{\mu} - \mu - \hat{\mu}^* + \mu^*)^T \Sigma_p^{-1}(\hat{\mu} - \mu^* + \mu^*) = O_p(1). \]

From Theorem 5.2 it follows that
\[ 2N(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\hat{\mu} - \mu - \hat{\mu}^* + \mu^*) = O_p(\sqrt{N}). \]

The last term we need to consider is
\[ N(1 + \zeta)^{-1}(\hat{\mu} - \mu - \hat{\mu}^* + \mu^*) . \]

Since \( \Sigma_p^{-1} \) is positive definite it follows that
\[ (\mu - \mu^*)^T \Sigma_p^{-1}(\mu - \mu^*) > 0, \]
and we have
\[ N(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\mu - \mu^*) \rightarrow \infty. \]

Furthermore when we divide the above by \( \sqrt{N} \) we get
\[ N^{1/2}(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\mu - \mu^*) \rightarrow \infty. \]

Therefore \( N(1 + \zeta)^{-1}(\mu - \mu^*)^T \Sigma_p^{-1}(\mu - \mu^*) \) dominates all the other terms in the limit and the theorem follows.

**Theorem 5.4** Suppose that Assumptions 2.1, 2.3, 2.4, (3.3), (3.5), and (3.12) hold. Then for each fixed \( p \geq 1 \) and \( r \geq 1 \), we have
\[ N^{1/2}(\hat{\mu}_v - \mu_v) \overset{d}{\rightarrow} N\left(0, \Sigma_e \otimes \Gamma^{-1} + E\left[\Delta_1 \otimes (\Gamma^{-1} \Delta_2 \Gamma^{-1})\right]\right) \]
where \( I_r \) is the \( r \times r \) identity matrix, and
\[ \Delta_1(j, t) = \left(\sum_{s=p+1}^{\infty} \mu_{sj} \langle v_s, X_1 \rangle\right) \left(\sum_{x=p+1}^{\infty} \mu_{xt} \langle v_x, X_1 \rangle\right), \]
and
\[ \Delta_2(i, q) = \langle v_i, X_1 \rangle \langle v_q, X_1 \rangle. \]

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Proof: The asymptotic normality follows from an application of the multivariate CLT. The derivation of the exact form of the asymptotic variance involves lengthy technical manipulations, and is omitted to conserve space.

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