

# Inference for Isolated Matrices and Structured Families of Matrices

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## Abstract

In this article we show how to use *vec* type operators to validate models for symmetric stochastic matrices. *vec* -type operators are operators that they also match vectors to matrices, that is, they allow to reorganize some elements of a matrix into a column vector. The presented formulation allowed us to make inference about the series of study, because the results presented in this work can be applied to the matrices of Hilbert-Schmidt products, that are very important matrices in the first phase of the STATIS methodology. Thus, the operators allowed us to present results that allow inference to be made for models for symmetric stochastic matrices. These models provided the basis for making inference for isolated matrices and structured families of matrices. In particular, we consider the case in which the matrices correspond to the treatments of base models. Thus, using the results presented, using *vec* type operators, it is possible to adjust a model with degree  $s$ . If  $\lambda_s = 0$ , the model can be simplified. These results, as already mentioned, can be applied to the matrices of Hilbert-Schmidt products (which are very important in the first phase of the STATIS Methodology), and to cross-product matrices which have an important role in inference.

## Keywords

Vec type operator's, Cross product matrices, Analysis of Variance, Models, Structured families.

## 1. Introduction

We use *vec* type operators in studying models for symmetric stochastic matrices based on their spectral decomposition. Namely we study  $k$ -degree models of the form

$$M = \mu + \bar{E} = \sum_{i=1}^k \lambda_i \alpha_i \alpha_i^t + \bar{E}$$

for symmetric stochastic matrices  $\mathbf{M}$ , with mean matrix  $\boldsymbol{\mu}$ , and  $\bar{\mathbf{E}}$  is a symmetric stochastic matrix with null mean. The models are developed using the spectral analysis of the matrices  $\boldsymbol{\mu}$ .

The adjustment and validation of the model requires the usage of the vector  $\tilde{\boldsymbol{\beta}}_i = \theta_i \boldsymbol{\gamma}_i, i = 1, \dots, k$ , which is an estimator of the structure vector  $\boldsymbol{\beta}_i = \lambda_i \boldsymbol{\alpha}_i, i = 1, \dots, k$  of  $\mathbf{M}$  with  $\lambda_1, \dots, \lambda_k$  the non-null eigenvalues and  $\boldsymbol{\alpha}_1, \dots, \boldsymbol{\alpha}_k$  are the eigenvectors, of the matrix  $\mathbf{M}$ .

We intend to introduce a new formulation for the stochastic models of symmetric matrices. In this new formulation, we will use  $\mathbf{vec}$  type operators to validate models for symmetric stochastic matrices. These operators, who play an important part in the new formulation for these models also facilitate the presentation of these results.

## 2. Model

The models are developed using the spectral analysis of the matrices  $\boldsymbol{\mu}$

$$\mathbf{M} = \boldsymbol{\mu} + \bar{\mathbf{E}} = \sum_{i=1}^k \lambda_i \boldsymbol{\alpha}_i \boldsymbol{\alpha}_i^t + \bar{\mathbf{E}},$$

where  $\bar{\mathbf{E}} = \frac{1}{2}(\mathbf{E} + \mathbf{E}^t)$  with  $\mathbf{vec}(\mathbf{E}) \sim N(\mathbf{0}, \sigma^2 \mathbf{I}_{n^2})$ , as we see in (Dias et al., 2013). The rank of the mean matrix will be the model's degree. The  $\mathbf{vec}$  type operators are operators are also vectors correspond to matrices. Namely for a symmetric matrix  $\mathbf{C} = [c_{ij}]$  of order  $k$ , we have

$$s\_vec(\mathbf{C}) = \begin{bmatrix} c_{11} \\ \vdots \\ c_{pp} \\ \vdots \\ c_{kk} \\ \sqrt{2}c_{12} \\ \vdots \\ \sqrt{2}c_{(k-1)k} \end{bmatrix}$$

$$w\_vec(\mathbf{C}; p) = \begin{bmatrix} c_{11} \\ \vdots \\ c_{pp} \\ \sqrt{2}c_{12} \\ \vdots \\ \sqrt{2}c_{pk} \end{bmatrix} \quad \text{and} \quad z\_vec(\mathbf{C}; p) = \begin{bmatrix} c_{(p+1)(p+1)} \\ \vdots \\ a_{kk} \\ \vdots \\ \sqrt{2}c_{(p+1)(p+2)} \\ \vdots \\ \sqrt{2}c_{(k-1)k} \end{bmatrix},$$

where

- $s\_vec(\mathbf{C})$  is a vector with  $g$  components where  $g = \frac{k(k+1)}{2}$ ;
- $w\_vec(\mathbf{C}; p)$  is a vector with  $c_p$  components where  $c_p = g - z_p$ ;
- $z\_vec(\mathbf{C}; p)$  is a vector with  $z_p$  components where  $z_p = \frac{(k-p)(k-p+1)}{2}$ .

We have also

$$\mathbf{v}_j(\mathbf{C}) = \begin{bmatrix} c_{jj} \\ c_{j(j+1)} \\ \vdots \\ c_{j(k-1)} \\ \vdots \\ c_{jk} \end{bmatrix}$$

which is a column vector whose components are the components of the  $j - th$  row vector of matrix  $\mathbf{C}$  starting at position  $(j, j)$ . The vector  $\mathbf{v}_j(\mathbf{C})$  have  $a_j$  components where  $a_j = k - j + 1$ , (Harville 1998), (Schott 1997), (Salvador, 2013).

### 3. Adjustment and validation of the model

Following Oliveira and Mexia (2007) we estimate the structure vector  $\beta_i = \lambda_i \alpha_i$ ,  $i = 1, \dots, k$  by the  $\tilde{\beta}_i = \theta_i \gamma_i$ ,  $i = 1, \dots, k$ , where  $\theta_1, \dots, \theta_n$  are the eigenvalues and the  $\gamma_1, \dots, \gamma_n$  the eigenvectors of matrix  $\mathbf{M}$ . With  $m_1, \dots, m_n$  the column vectors of  $\mathbf{M}$ , we have

$$\tilde{\beta}_i = \mathbf{M}\gamma_i = \mathbf{M}^t \gamma_i = \begin{bmatrix} m_1^t \gamma_i \\ \vdots \\ m_n^t \gamma_i \end{bmatrix} = \begin{bmatrix} \gamma_i^t m_1 \\ \vdots \\ \gamma_i^t m_n \end{bmatrix} = (\mathbf{I}_n \otimes \gamma_i^t) \begin{bmatrix} m_1 \\ \vdots \\ m_n \end{bmatrix} \mathbf{Z}, i = 1, \dots, k,$$

so  $\tilde{\beta}_i = \theta_i \gamma_i$ ,  $i = 1, \dots, k$  will have mean vector and variance-covariance matrix

$$\begin{cases} E(\tilde{\beta}_i) = (\mathbf{I}_n \otimes \gamma_i^t) \boldsymbol{\eta}, i = 1, \dots, k \\ \Sigma(\tilde{\beta}_i) = \sigma^2 (\mathbf{I}_n \otimes \gamma_i^t) \mathbf{L} (\mathbf{I}_n \otimes \gamma_i), i = 1, \dots, k \end{cases}$$

If  $\mathbf{M}$  has degree  $k$ , and the  $\tilde{\beta}_1, \dots, \tilde{\beta}_k$  [ $\gamma_1 \dots \gamma_k$ ] are good estimators of  $\beta_1, \dots, \beta_k$  [ $\alpha_1 \dots \alpha_k$ ] then, we can write

$$\bar{\mathbf{E}} = \mathbf{M} - \sum_{i=1}^k \beta_i \alpha_i^t \approx \mathbf{M} - \sum_{i=1}^k \tilde{\beta}_i \gamma_i^t = \mathbf{M} - \sum_{i=1}^k (\mathbf{I}_n \otimes \gamma_i^t) \mathbf{Z} \gamma_i^t$$

and then

$$\mathbf{R} = \text{vec}(\bar{\mathbf{E}}) \approx \mathbf{Z} - \text{vec} \left( \sum_{i=1}^k ((\mathbf{I}_n \otimes \gamma_i^t) \mathbf{Z}) \gamma_i^t \right),$$

with

$$\text{vec} \left( \sum_{i=1}^k ((\mathbf{I}_n \otimes \gamma_i^t) \mathbf{Z}) \gamma_i^t \right) = \sum_{i=1}^k (\gamma_i \otimes \mathbf{I}_n \otimes \gamma_i^t) \mathbf{Z}.$$

Thus

$$\mathbf{R} = (\mathbf{I}_{n^2} - \mathbf{W}) \mathbf{Z},$$

with  $\mathbf{W} = \sum_{i=1}^k (\gamma_i \otimes \mathbf{I}_n \otimes \gamma_i^t)$  and  $\Sigma(\mathbf{R}) = \sigma^2 (\mathbf{I}_{n^2} - \mathbf{W}) \mathbf{L} (\mathbf{I}_{n^2} - \mathbf{W})^t$ .

While the vector  $\mathbf{R}$  can be considered as a residual vector,  $\tilde{\beta}_1, \dots, \tilde{\beta}_k$  [ $\gamma_1 \dots \gamma_k$ ] will be the adjusted global structure vector.

Now to check if the model holds it is possible to derive from  $\mathbf{R}$  a vector that, when the model holds, is normal and homoscedastic, (Areia et al. 2008). We then apply normality tests to that vector.

#### 4. Application to the Inter-structure

In the Joint Table Analysis (ACT), it is customary to have several double-entry data tables, the data that make up these tables are quantitative in nature, which can be obtained in different situations. Each table consists of  $n$  rows (objects) and  $p$  columns (variables). The various techniques developed for the analysis of this type of tables date back to the 70s and 80s, namely studies of these techniques in Escoufier and L'Hermier des Plantes (1976), Foucart (1981), among others. The various techniques developed allow for a global comparison of all tables, as well as the study of the possible existence of a common structure, which is called inter-structure. The inter-structure is the first stage of the STATIS methodology (STATIS and STATIS Dual method). This methodology was introduced by Hermier des Plantes (1976) and was developed in particular by Lavit (1988). The STATIS methodology uses Euclidean distances between configurations observed in  $k$  different situations that we call studies. Whenever, the variables differ throughout the  $k$  studies and the number of objects is fixed, there is a proximity between the individuals (objects) we say that we are within the scope of the STATIS method. Whenever the reverse occurs, that is, the number of variables is fixed and objects can change over the course of studies, we say that we are within the scope of the STATIS Dual method. In these two methods, series of studies are studied in which a trio of matrices is considered  $(\mathbf{X}, \mathbf{D}, \dot{\mathbf{D}})$ , constituted by a data matrix  $\mathbf{X}$  and two weight matrices, one for objects and one for variables,  $\mathbf{D}$  and  $\dot{\mathbf{D}}$  respectively. When the objects are the same for all studies in a series, the same is said of the first type. When the variables are the same for all studies in a series, the same is said for the second type. Escoufier (1978) defined the operators

$$\mathbf{A}_i = \mathbf{X}_i \dot{\mathbf{D}}_i, \mathbf{X}_i^t \mathbf{D}_i, i = 1, \dots, n,$$

and

$$\mathbf{B}_i = \mathbf{X}_i^t \mathbf{D}_i \mathbf{X}_i \dot{\mathbf{D}}_i, i = 1, \dots, n,$$

to represent the studies

$$(\mathbf{X}_i, \mathbf{D}_i, \dot{\mathbf{D}}_i), i = 1, \dots, n,$$

in the case of series of first and second types, respectively. Let's consider the study of the first type series, putting  $\mathbf{U} = \mathbf{A}$ , then we have the matrices  $\mathbf{S} = [s_{ij}]$ , with

$$s_{ij} = \text{tr}(\mathbf{U}_i \mathbf{U}_j^t), i = 1, \dots, n,$$

the product of Hilbert-Schmidt operators  $\mathbf{U}_i \in \mathbf{U}_j, i = 1, \dots, n$ . Being  $\theta_1 \geq \dots \geq \theta_n$ , the eigenvalues and  $\boldsymbol{\gamma}_1, \dots, \boldsymbol{\gamma}_n$  the corresponding eigenvectors, and according Escoufier (1978), the  $j$ -th study is represented by the point whose coordinates are the  $j$ -th components of the vectors  $\theta_1^{\frac{1}{2}} \boldsymbol{\gamma}_1, \dots, \theta_n^{\frac{1}{2}} \boldsymbol{\gamma}_n$ . When these points are along the first axis with the direction of  $\boldsymbol{\gamma}_1, \theta_1$  it prevails over the other eigenvalues and the series has one-degree structure (Lavit, 1988). In that case it can be admitted having

$$\mathbf{S} = \lambda \boldsymbol{\alpha} \boldsymbol{\alpha}^t + \bar{\mathbf{E}},$$

therefore, having a symmetric stochastic matrix with degree 1. Thus, it can be admitted that

$$\mathbf{U}_i = \mathbf{U}_i^0 + \mathbf{E}_i, i = 1, \dots, n,$$

com  $\mathbf{U}_i^0 = [u_{ilh}]$  and  $\mathbf{E}_i = [e_{ilh}]$  with  $\text{vec}(\mathbf{E}_i) \sim N(0, \sigma^2 \mathbf{I}_{m^2}), i = 1, \dots, n$  (Arcia, 2009).

It is also admitted that  $\text{vec}(\mathbf{E}_i)$  are independent. Thus,

$$s_{ij} = \text{tr}(\mathbf{U}_i \mathbf{U}_j^t) = \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} + e_{ilh})(\mathbf{u}_{jlh} + e_{jlh})$$

$$\begin{aligned}
 &= \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} \mathbf{u}_{jlh} + \mathbf{u}_{ilh} e_{jth} + \mathbf{u}_{jth} e_{ilh} + e_{ilh} e_{jth}) \\
 &= S_{ij}^0 + \sum_{l=1}^m \sum_{h=1}^m (e_{ilh} e_{jth}), \quad i, j = 1, \dots, n,
 \end{aligned}$$

with  $S_{ij}^0 = \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} \mathbf{u}_{jlh} + \mathbf{u}_{ilh} e_{jth} + \mathbf{u}_{jth} e_{ilh})$ ,  $i, j = 1, \dots, n$ .

With  $\mathbf{S}^0 = [S_{ij}^0]$ ,  $\mathbf{vec}(\mathbf{S}^0)$  it will be normally distributed. We can therefore admit that  $\mathbf{Z} = \mathbf{vec}(\mathbf{S})$  it will be (approximately) normally distributed as long as we can despised  $\sum_{l=1}^m \sum_{h=1}^m (e_{ilh} e_{jth})$ ,  $i, j = 1, \dots, n$ . Through simulations carried out see (Dias, 2013, p.85-95), it is shown that when the coefficients of variation  $\frac{\sigma_j}{\mu_j}$ , of a pair of normal variables  $X_j, j = 1, 2$  are less than  $\frac{1}{17}$ , the product  $X_1 X_2$  can be treated as being normally distributed. So, when the matrix elements  $\mathbf{U}_1, \dots, \mathbf{U}_n$  have coefficients of variation less than  $\frac{1}{17}$  we can treat  $\mathbf{Z}$  as normal.

With

$$\mathbf{K}_{ij} = \text{tr}(\mathbf{U}_i^0 \mathbf{U}_j^{0t}) = \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} + \mathbf{u}_{jlh}), \quad i, j = 1, \dots, n,$$

given the independence and homoscedasticity of  $\mathbf{vec}(\mathbf{E}_i)$ ,  $i = 1, \dots, n$ , we have

$$\begin{aligned}
 \text{Var}(S_{ij}^0) &= \text{Var} \left( \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} + e_{ilh} + \mathbf{u}_{jth} + e_{jth}) \right) = \sigma^2 \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh}^2 + \mathbf{u}_{jth}^2) \\
 &= \sigma^2 (\mathbf{K}_{ii} + \mathbf{K}_{jj}), \quad i \neq j,
 \end{aligned}$$

as well as

$$\begin{aligned}
 \text{Var}(S_{ij}^0) &= \text{Var} \left( 2 \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh} + e_{ilh}) \right) = 4\sigma^2 \sum_{l=1}^m \sum_{h=1}^m (\mathbf{u}_{ilh}^2) \\
 &= 4\sigma^2 (\mathbf{K}_{ii}), \quad i = 1, \dots, n.
 \end{aligned}$$

The remaining elements of the variance-covariance matrix ( $\mathbf{L}$ ) are (approximately) given by (Areia, 2009),

$$\begin{cases}
 \text{Cov}(S_{ii}, S_{ij}) = \text{Cov}(S_{ii}, S_{ji}) = 2\sigma^2 \mathbf{K}_{ij}, \quad i \neq j. \\
 \text{Cov}(S_{ii}, S_{ji}) = 0 \quad i \neq j \\
 \text{Cov}(S_{ij}, S_{ji}) = \text{Var}(S_{ij}) \sigma^2 (\mathbf{K}_{ii} + \mathbf{K}_{jj}), \quad i \neq j \\
 \text{Cov}(S_{ij}, S_{il}) = \text{Cov}(S_{ij}, S_{li}) \approx \sigma^2 \mathbf{K}_{jl}, \quad i \neq j \neq l \\
 \text{Cov}(S_{ij}, S_{lu}) = 0, \quad i \neq j \neq l \neq u
 \end{cases}$$

We can then admit (with due approximation) that  $\mathbf{Z} \sim N(\boldsymbol{\eta}, \sigma^2 \mathbf{L})$ , with  $\boldsymbol{\eta} = \mathbf{vec}([\mathbf{K}])$  and  $\mathbf{K} = [k_{ij}]$ , since the components of  $\mathbf{vec}(\mathbf{S}^0)$  being linear functions of a normal vector will have normal joint distribution. Since the  $\mathbf{S}$  model is a particular case of the base model

$$\mathbf{M} = \boldsymbol{\mu} + \bar{\mathbf{E}} = \sum_{i=1}^k \lambda_i \boldsymbol{\alpha}_i \boldsymbol{\alpha}_i^t + \bar{\mathbf{E}}$$

with  $(\lambda_i, \boldsymbol{\alpha}_i)$  are the pairs (eigenvalues, eigenvectors) of  $\boldsymbol{\mu}$  and  $\bar{\mathbf{E}}$  is a symmetric stochastic matrix. with  $(\lambda_i, \boldsymbol{\alpha}_i)$  are the pairs (eigenvalues, eigenvectors) of  $\boldsymbol{\mu}$  and  $\bar{\mathbf{E}}$  is symmetric stochastic matrix. To avoid considering a possible large

number of small eigenvalues we assume that  $\lambda_1 \geq \dots \geq \lambda_k > c$ . With  $\beta_i = \lambda_i \alpha_i, i = 1, \dots, k$  the information contained in  $\mathbf{M}$  can be condensed in a structured vector

$$\boldsymbol{\beta} = [\boldsymbol{\beta}_1^t \dots \boldsymbol{\beta}_k^t]^t,$$

and, a sum of square of residues

$$V = \|\mathbf{M}\|^2 - \|\boldsymbol{\beta}\|^2,$$

where  $\|\cdot\|$  indicates Euclidean norm both for matrices and vectors.

In this article we use the *vec* and other relater operators to carry out inference for structured families of symmetric stochastic matrices  $\mathbf{M}$ . These are obtained through the sum of the respective mean matrix and a symmetric stochastic matrix with null mean. We consider that the *vec* operator of the matrix  $\mathbf{E}$  is normal homoscedastic. The matrices on these families correspond to the treatments of a base design, and the inference is centered on model validation and the action of the factors in the base model on mean matrices.

### 3. Application to Cross Product Matrices

Given the data matrix  $\mathbf{X}$  of type  $v \times d$ , the matrices  $\mathbf{X}\mathbf{X}^t$  and  $\mathbf{X}^t\mathbf{X}$  are the matrices of cross products for row and column vector of  $\mathbf{X}^t\mathbf{X}$ . In particular if the column vectors of  $\mathbf{X}$  are the *vec*( $\mathbf{U}_i$ ),  $i = 1, \dots, d$ ,  $\mathbf{X}^t\mathbf{X}$  will be the matrix of Hilbert-Schmidt products of the matrices  $\mathbf{U}_1, \dots, \mathbf{U}_d$ . Thus, these models directly generalize the previous ones, making it easy to verify that the treatment presented extends directly.

In the case of the matrix  $\mathbf{X}\mathbf{X}^t$  we should have

$$k_{ij} = \sum_{l=1}^d x_{il}x_{jl}, i, j = 1, \dots, m,$$

with  $\mathbf{X} = [x_{ih}]$ . For the matrix  $\mathbf{X}^t\mathbf{X}$  we will have

$$k_{ij} = \sum_{l=1}^m x_{li}x_{lj}, i, j = 1, \dots, d.$$

Note that the matrices  $\mathbf{X}\mathbf{X}^t$  and  $\mathbf{X}^t\mathbf{X}$  have the same non-null eigenvalues  $\theta_1 \geq \dots \geq \theta_r > 0$ , with  $r = \text{Car}(\mathbf{X}) = \text{Car}(\mathbf{X}\mathbf{X}^t) = \text{Car}(\mathbf{X}^t\mathbf{X})$ . Thus, the models for  $\mathbf{X}\mathbf{X}^t$  and  $\mathbf{X}^t\mathbf{X}$  will have the same degree  $k'$ . Assuming that  $d < v$ , the model must be adjusted to  $\mathbf{X}^t\mathbf{X}$  which will be a matrix  $d \times d$  rather than  $\mathbf{X}\mathbf{X}^t$  that is  $v \times v$ . At  $\boldsymbol{\gamma}_1, \dots, \boldsymbol{\gamma}_{k'}$  first  $k'$  eigenvectors for  $\mathbf{X}^t\mathbf{X}$  will correspond, as eigenvectors of  $\mathbf{X}\mathbf{X}^t$ , the

$$\boldsymbol{\gamma}_i = \frac{1}{\sqrt{\theta_i}} \mathbf{X}\boldsymbol{\gamma}_i, i = 1, \dots, k',$$

that can be used in the inference for the  $\mathbf{X}\mathbf{X}^t$  models and their families.

Since  $V = \|\mathbf{Y}\|^2$ , with  $V$  the sum of the squared residuals, and considering  $\mathbf{Y}$  normal with mean vector and covariance matrix  $\sigma^2 \mathbf{I}_m$ , is independent from the structure vector  $\dot{\boldsymbol{\beta}}_i$ , with

$$\dot{\boldsymbol{\beta}}_i = (\mathbf{X}^t\mathbf{X})\boldsymbol{\gamma}_i, i = 1, \dots, k'$$

that is independent of

$$\boldsymbol{\beta}_i = \mathbf{X}\dot{\boldsymbol{\beta}}_i, i = 1, \dots, k',$$

so, the results for the model corresponding to  $\mathbf{X}^t\mathbf{X}$  are carried directly to the model corresponding to  $\mathbf{X}\mathbf{X}^t$ .

In many cases grade one models can be used for these matrices (Oliveira 2001), (Ramos 2007) and (Areia 2011).

#### 4. Structured Families

These vectorial operators, besides presenting themselves as an important part in the new formulation for these models, also facilitate the presentation of the results. The models for symmetric stochastic matrices are the basis for inference for isolated matrices and for structured families of matrices (Dias 2013), (Dias et al. 2016).

In these families the matrices, all of the same order, correspond to the treatments of base models. Since the matrices have all the same order, we are in the balanced case where we have the same number of degrees of freedom for the error for each treatment.

The ANOVA and related techniques are, in the balanced case, robust techniques for heteroscedasticity and even more for non-normality (Ito 1980), (Scheffé 1959).

The symmetric  $n \times n$  matrix  $W = [w_{ij}]$  has, besides its *vec* the *semi\_vecS*( $M$ ) with components  $w_{1,1}, \dots, w_{n,n}, w_{2,3}, \dots, w_{2,n}, \dots, w_{n-1,n}$ .

We use the *semi\_vec* to validate models for symmetric stochastic matrices. When the predominance of the first eigenvalue is very large, we can adopt a model of degree one of the form

$$M = \mu + \bar{E},$$

where  $\mu$  is a mean matrix and  $\bar{E}$  is a symmetric stochastic matrix with null mean, such that  $\bar{E} = \frac{1}{2}(E + E^t)$  with  $\text{vec}(E) \sim N(0, \sigma^2 I_{n^2})$  normal homoscedastic.

These models, as we mentioned earlier, have been used successfully in several applications, namely the first phase of the STATIS methodology, the inter-structure. Since Hilbert-Schmidt matrices are matrices of cross products, we can use them for the latter, changing to higher-grade models if degree one models do not fit. In what follows we will base ourselves on presenting the main results that we have for these models.

Simulations (presented next, in the Table 1), show that when the first eigenvalue is sufficiently dominant, we can conclude that the mean matrix as rank one and that the first eigenvalue and eigenvector can be used to estimate the sole non null eigenvalue of the mean matrix and the corresponding eigenvector.

It is intended to generalize this approach, that is, when the first  $n$  eigenvalues of the matrix  $M$  are sufficiently larger than the others, we can use them to estimate  $\lambda_1, \dots, \lambda_r$  and the corresponding eigenvectors to estimate  $\alpha_1, \dots, \alpha_r$ .

Thus, with  $\theta_1 > \theta_2 > \dots > \theta_n$  the eigenvalues of the matrix  $M$  associated to the eigenvectors  $\gamma_1, \dots, \gamma_r$ , respectively. Thus

$$M = \sum_{i=1}^n \theta_i \gamma_i \gamma_i^t.$$

The first  $r$  eigenvectors of structure can be estimated by  $\tilde{\beta}_i = \theta_i \gamma_i$ , with  $i = 1, \dots, n$ .

The most interesting cases for these types of models occur when the matrices of this family have the first dominant eigenvalue or the first dominant eigenvalues. In that case, the action of the factors of that design, on the components of the first structure vector or the first structure vectors, depending as it is only the first dominant eigenvalue or the first dominant eigenvalues, respectively (Dias et al. 2012), (Salvador 2013).

Let,  $(\theta_i, \gamma_i)$   $i = 1, \dots, n$  be the pairs of eigenvalues and eigenvectors for  $M$ . Then

$$\|M\|^2 = \sum_{i=1}^n \theta_i^2,$$

when

$$R = \frac{\theta_i^2}{\sum_{l=i+1}^k \theta_l^2}$$

If  $c_i$  is large we can take the pair  $(\theta_i, \gamma_i)$   $i = 1, \dots, n$  as estimators of  $(\lambda_i, \alpha_i)$   $i = 1, \dots, n$ , (Oliveira and Mexia 2004). The values considered for  $k$  were: 2, 6, 8, 14, 17 and 40 and the values of  $R$  were 12, 5; 50, 100, 200, 600 and 800.

In the construction of the  $M$  matrices we took advantage of the symmetry of the problem taking  $\alpha = \zeta_1$ .

For each pair  $(k, R)$ , 1000 matrices  $\mathbf{E}$  were generated with

$$\text{vec}(\mathbf{E}) \sim N(\mathbf{0}, \mathbf{I}_{k^2})$$

from which we obtained the matrices

$$\mathbf{W} = \lambda \boldsymbol{\zeta}_1 \boldsymbol{\zeta}_1^t + \bar{\mathbf{E}},$$

with

$$\lambda = \sqrt{2R},$$

and  $\boldsymbol{\alpha} = \boldsymbol{\zeta}_1$ . Being  $\boldsymbol{\gamma}_1$  the first eigenvector of  $\mathbf{W}$ , since

$$\|\boldsymbol{\zeta}_1\|^2 = \|\boldsymbol{\gamma}_1\|^2 = 1,$$

we will have

$$\begin{aligned} \|\boldsymbol{\zeta}_1 - \boldsymbol{\gamma}_1\|^2 &= (\boldsymbol{\zeta}_1 - \boldsymbol{\gamma}_1)^t (\boldsymbol{\zeta}_1 - \boldsymbol{\gamma}_1) \\ &= \boldsymbol{\zeta}_1^t \boldsymbol{\zeta}_1 - 2\boldsymbol{\zeta}_1^t \boldsymbol{\gamma}_1 + \boldsymbol{\gamma}_1^t \boldsymbol{\gamma}_1 \\ &= 2(1 - \boldsymbol{\gamma}_{11}), \end{aligned}$$

since  $\boldsymbol{\gamma}_{11} = \boldsymbol{\zeta}_1^t \boldsymbol{\gamma}_1$ , with  $\boldsymbol{\gamma}_{11}$  the first component of  $\boldsymbol{\gamma}_1$ . If  $\|\boldsymbol{\gamma}_1\|^2 = 1$ , we will have  $-1 \leq \boldsymbol{\gamma}_{11} \leq 1$ , and

$$\|\boldsymbol{\alpha} - \boldsymbol{\gamma}_1\|^2 = \|\boldsymbol{\zeta}_1 - \boldsymbol{\gamma}_1\|^2,$$

is much smaller as much closer to 1 is  $\boldsymbol{\gamma}_{11}$ . Thus, we can consider  $\boldsymbol{\gamma}_{11}$  as a measure of efficiency of  $\boldsymbol{\gamma}$ , to estimate  $\boldsymbol{\alpha}$ .

Thus, for each pair  $(k, R)$ , we present in Table 1 results for batches of 1000 simulations, and the respective mean values and standard deviation for  $\|\boldsymbol{\alpha} - \boldsymbol{\gamma}_1\|^2$ . From the analysis of the Table 1 we see that, there are many times when the estimated values for  $c_i$  fall in the zone where we are led to admit having  $\boldsymbol{\gamma} \approx \boldsymbol{\alpha}$  and consequently,  $\theta_k \approx \lambda$ , and when  $R > 200$ , we can use  $\theta$  and  $\boldsymbol{\gamma}$  as estimators of  $\lambda$  and  $\boldsymbol{\alpha}$  respectively.

Table 1. Mean values and standard deviations for  $\|\boldsymbol{\alpha} - \boldsymbol{\gamma}_1\|^2$ .

R	k = 2		k = 6		k = 8	
	Mean	S.D.	Mean	S.D.	Mean	S.D.
12.5	0.944	0.181	0.691	0.288	0.501	0.287
50	0.941	0.061	0.908	0.118	0.878	0.128
100	0.986	0.008	0.973	0.010	0.965	0.011
200	0.998	0.002	0.997	0.002	0.996	0.003
600	0,999	0,000	0,999	0,001	0,998	0,001
800	0,999	0,000	0,999	0,000	0,000	0,000
R	k = 14		k = 17		k = 40	
	Mean	S.D.	Mean	S.D.	Mean	S.D.
12.5	0,432	0.296	0,378	0.263	0,417	0.243
50	0,787	0.165	0,891	0.177	0,705	0.223

100	0,943	0.038	0,932	0.051	0,869	0.062
200	0,996	0.012	0,956	0.016	0,889	0.016
600	0,998	0,002	0,997	0,004	0,983	0,002
800	0,999	0,000	0,999	0,000	0,999	0,000

Let us put

$$\bar{M}_h = M - \sum_{i=1}^h \theta_i \gamma_i \gamma_i^t,$$

where  $m_h = S(\bar{M}_h)$ . If  $k \leq h$  we will have  $\bar{M}_h \approx E$ , and so the components of  $S(\bar{M}_h)$  can be assumed to be normal with null mean values.

With  $Y_1, \dots, Y_g$  ( $g = n(n-1)$ ) the components of  $S(\bar{M}_h)$  we are then led to test

$$H_{0k}: Y_1, \dots, Y_g \text{ i. i. d. } \sim N(0, \sigma^2),$$

using the statistics (Mexia 1989, 1989b)

$$F = \frac{(g-1)(\sum_{i=1}^g Y_i)^2}{g \sum_{i=1}^g Y_i^2 - (\sum_{i=1}^g Y_i)^2},$$

which when the hypothesis holds, has the central F distribution with 1 and  $g-1$  degrees of freedom (Mexia 1995). When  $H_0$  thus not hold, the  $Y_1, \dots, Y_g$  will have mean values  $\mu_1, \dots, \mu_g$  and the numerator and denominator of F will have the non-centrality parameter

$$\delta_1 = \frac{1}{\sigma^2} \left( \frac{1}{g} \sum_{i=1}^g \mu_i \right)^2, \text{ and } \delta_2 = \frac{1}{\sigma^2} \left( \sum_{i=1}^g \mu_i^2 - (\sum_{i=1}^g Y_i)^2 \right),$$

hence, there will be alternatives in which  $\delta_1$  predominates over  $\delta_2$  ( $\delta_2$  predominates over  $\delta_1$  and in which  $F$  tends to take larger (smaller) values than when  $H_0$  holds. A numerical study showed that, with  $f_{p,1,g-1}$  the quantile for probability  $p$  of the central F distribution with 1 and  $g-1$  degrees of freedom, the acceptance region.

$$\left[ f_{\frac{p}{2},1,g-1}; f_{1-\frac{p}{2},1,g-1} \right],$$

leads to, p level tests with power increasing rapidly with  $\delta_1$  and  $\delta_2$  (Mexia 1995).

We test this hypothesis for increasing values of k. With  $\bar{k}$  the smallest value for which the hypothesis is not rejected we will have the degree of the adjusted model given by

$$M^* = \sum_{i=1}^{\bar{k}} \theta_i \gamma_i \gamma_i^t.$$

We can now complete the adjustment testing hypothesis that  $\gamma_{\bar{k}+1}, \dots, \gamma_n, \bar{k} = 1, \dots, n-1$  are normal homoscedastic with null mean vectors independent between them self's, and from  $\gamma_{\bar{k}}$ , which will have a mean vector  $\gamma_{\bar{k}} \neq 0$ , using the test statistic

$$\bar{F}_{\bar{k}} = (n - \bar{k}) \frac{\|\gamma_{\bar{k}}\|^2}{\sum_{h=\bar{k}+1}^n \|\gamma_h\|^2}.$$

This F test has  $n$  and  $(n - \bar{k})n$  degrees of freedom and non-centrality parameter

$$\delta_h = \frac{1}{\sigma^2} \|\mathbf{r}_k\|^2,$$

and the p level critical value  $F_{1-p,n,(n-k)n}$ .

## 5. Conclusion

In this article we show how to use *vec* type operators to validate the model. The presented formulation allowed us to make inference about the series of study, because the results presented in this work can be applied to the matrices of the products of Hilbert-Schmidt and cross product matrices. These operators allowed us to present results that allow inference to be made for models for symmetric stochastic matrices. The *semi\_vec* can be used to determinate the number of non-null eigenvalues of a symmetric matrix  $\mathbf{M} = \boldsymbol{\mu} + \bar{\mathbf{E}}$ , and we obtained, using adjusted eigenvalues and eigenvectors, an adjusted model for  $\mathbf{M}$ . These models provided the basis for making inference for isolated matrices and structured families of matrices.

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