

# Sparse Online Principal Component for Laplace Factor Models

Siqi Liu, and Guangbao Guo

**Abstract**—This paper proposes a Sparse Online Principal Component (SOPC) approach applied to the Laplace Factor Model (LFM). The method integrates the strengths of Sparse Principal Component (SPC) and Online Principal Component (OPC) techniques, enabling real-time data updates and yielding sparse solutions. Numerical simulations validate its stability and accuracy under varying sample sizes and dimensions. Results demonstrate low error rates and robust adaptability in Laplace factor models.

**Index Terms**—Laplace factor model; sparse online principal component; stability and sensitivity; real-time updating; numerical simulation.

## I. INTRODUCTION

THE Laplace Factor Model (LFM) is widely used for analyzing high-dimensional multivariate data by extracting latent factors. Existing methods face limitations in real-time updating and sparsity. This study introduces the Sparse Online Principal Component (SOPC) method, which supports online updates and sparse parameter estimation. Simulations assess its sensitivity to sample size and dimensionality. The findings offer a practical tool for dynamic data analysis in complex scenarios.

## II. LAPLACE FACTOR MODEL

### A. Laplace distribution

The Laplace distribution encompasses three primary forms, defined by their probability density functions.

Consider a Laplace-distributed random variable  $X$ , parameterized by its location parameter  $\mu$ , scale parameter  $b$ , and precision parameter  $\tau = \frac{1}{b}$ .

1) *Univariate symmetric Laplace distribution:*

$$f_X(x; 0, b) = \frac{1}{2b} \exp\left(-\frac{|x|}{b}\right).$$

2) *Precision parameterization Laplace distribution:*

$$f_X(x; \mu, \tau) = \frac{\tau}{2} \exp(-\tau|x - \mu|).$$

3) *Univariate asymmetric Laplace distribution:*

$$f_x(x; \mu, b_1, b_2) = \begin{cases} \frac{1}{b_1} \exp\left(-\frac{|x-\mu|}{b_1}\right), & x < \mu, \\ \frac{1}{b_2} \exp\left(-\frac{|x-\mu|}{b_2}\right), & x \geq \mu. \end{cases}$$

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### B. Laplace regression

We consider a distributed Laplace regression model ( $k = 1, \dots, K_n$ ):

$$Y_{Ik} = X_{Ik}\gamma + \epsilon_{Ik}, \quad \epsilon_{Ik} \sim \text{Laplace}(\mu_1, b_3 I_{n_{ik} \times n_{ik}})$$

where

- $X_{Ik}$  is a submatrix of the full design matrix  $N_{Ik}$ , with  $n_{Ik} \geq p$ ;
- $\epsilon_{Ik}$  is a subresidual vector, and  $I_{n_k \times n_k}$  denotes the identity matrix;
- $\gamma = (\gamma_1, \dots, \gamma_p)^T$  denotes the regression coefficient vector;
- $\mu_1$  and  $b_3 > 0$  represent location and scale parameters, respectively.

We notice that

$$Y = (Y_{I_1}^T, Y_{I_2}^T, \dots, Y_{I_{K_n}}^T)^T,$$

$$X = (X_{I_1}^T, X_{I_2}^T, \dots, X_{I_{K_n}}^T)^T.$$

The model can also be expressed as:

$$Y = X\gamma + \epsilon, \quad \epsilon \sim \text{Laplace}(\mu_1, b_3),$$

with  $\mathbb{E}(Y) = \mu_1$ ,  $\text{Var}(\epsilon_{Ik}) = 2b_3^2 - \mu_1^2$  and the error variance estimator  $\hat{\sigma}^2 = \frac{\epsilon^T \epsilon}{n-p}$ .

### C. Laplace factor model

The formulation of the LFM is given by

$$X = FA^T + \epsilon, \quad \epsilon \sim \text{Laplace}(\mu_2, b_4 I_{p \times p}) \quad (1)$$

where

- $X \in \mathbb{R}^{n \times p}$  denotes the observed data matrix;
- $A \in \mathbb{R}^{p \times m}$  constitutes the factor loading matrix;
- $F \in \mathbb{R}^{n \times m}$  represents the matrix containing latent common factor with  $m < p$ ;
- $\epsilon \in \mathbb{R}^{n \times p}$  denotes the error matrix;
- $\mu_2$  and  $b_4$  represent location and scale parameters.

The assumptions of the model are as follows.

- The expected value of  $f_j$  is zero and its variance is  $I_{m \times m}$ .
- The expected value of  $\epsilon_j$  is zero and its variance is a diagonal matrix  $D = \text{diag}(\sigma_1^2, \sigma_2^2, \dots, \sigma_p^2)$ .
- The  $f_j$  and  $\epsilon_j$  are independent,  $\text{cov}(f_j, \epsilon_j) = 0$ .

## III. METHOD

### A. Sparse principal component analysis

To describe the covariance structure among  $p$  correlated variables, we use  $m \ll p$  common factors:

$$\Sigma = AA^T + D$$

where  $A$  denotes the factor loading matrix, and the diagonal matrix  $D$  accounts for specific variance.

Under the Sparse Principal Component (SPC) framework, eigendecomposition of the empirical covariance matrix  $S$  yields

$$S = \sum_{i=1}^p \lambda_i v_i v_i^T$$

where  $\lambda_i$  and  $v_i$  denote eigenvalues and their corresponding orthogonal eigenvectors. Let  $\Lambda = \text{diag}(\lambda_1, \lambda_2, \dots, \lambda_p)$  and  $V = (v_1, v_2, \dots, v_p)$ . Through sparsification, a sparse eigenvector matrix  $V_S$  is obtained, which is then used to derive the sparse loading matrix  $A_S$ .

Let  $V_s$  be initialized as  $V$ . The sparse optimization criterion is formulated as

$$V_S = \arg \min_{V_S} \left( \sum_{i=1}^n \|X_{i.} - W_S V_S^T X_{i.}\|_F^2 + \rho \sum_{j=1}^m \|V_{S,j}\|_F^2 + \theta \sum_{j=1}^m \|V_{S,j}\|_1 \right)$$

where  $W_S^T W_S = I_{m \times m}$  and  $n > p$  when  $\rho = 0$ . The sparsity parameter  $\theta > 0$  controls eigenvector sparsity. The matrix  $\hat{A}_S$  containing sparse loading and the matrix  $\hat{D}_S$  containing specific variance are calculated for  $i = 1, 2, \dots, p$ :

$$\hat{A}_S = (\sqrt{\lambda_1} v_{S1}, \sqrt{\lambda_2} v_{S2}, \dots, \sqrt{\lambda_m} v_{Sm}) = (\hat{a}_{Sij})_{p \times m},$$

$$\hat{D}_S = \text{diag}(\hat{\sigma}_1^2, \hat{\sigma}_2^2, \dots, \hat{\sigma}_p^2), \quad \hat{\sigma}_i^2 = s_{ii} - \sum_{j=1}^m \hat{a}_{Sij}^2.$$

### B. Sparse online principal component analysis

The SOPC method introduces a sparsity parameter  $\theta$  to enable online sparse eigenvector updates. For the first  $k$  observations  $X^k = (X_1^T, X_2^T, \dots, X_k^T)^T$ , ( $k < n$ ), the sample covariance matrix is  $S^k = V^k \Lambda^k (V^k)^T$ , with  $\Lambda^k = \text{diag}(\lambda_1^k, \lambda_2^k, \dots, \lambda_p^k)$ , and  $V^k = (v_1^k, v_2^k, \dots, v_p^k)$ . Upon receiving the  $(k+1)$ -th sample  $X_{k+1}$ , the covariance updates as

$$S^{k+1} = \frac{k}{k+1} S^k + \frac{1}{k+1} X_{k+1}^T X_{k+1}$$

and the eigen decomposition  $X_{k+1} = V^{k+1} \Lambda^{k+1} (V^{k+1})^T$  is performed. Let  $W_{SO} = V^{k+1} [1 : m]$ , and initialize  $V_{SO}^{k+1} = V^{k+1}$ . The sparse eigenvector matrix  $V_{SO}^{k+1}$  is obtained via

$$V_{SO}^{k+1} = \arg \min_{V_{SO}^{k+1}} \left( \|X_{k+1.} - W_{SO} V_{SO}^{k+1T} X_{k+1.}\|_F^2 + \rho \sum_{j=1}^m \|V_{SO,j}^{k+1}\|_F^2 + \theta \sum_{j=1}^m \|V_{SO,j}^{k+1}\|_1 \right),$$

where  $W_{SO}^T W_{SO} = I_{m \times m}$ . The optimization objective can be re-expressed as

$$L(V_{SO}^{k+1}) = \text{trace}(X_{k+1.}^T X_{k+1.}) + \sum_{j=1}^m [V_{SO,j}^{k+1T} (X_{k+1.}^T X_{k+1.} + \rho) V_{SO,j}^{k+1} - 2w_{SO,j}^T X_{k+1.}^T X_{k+1.} V_{SO,j}^{k+1} + \theta \|V_{SO,j}^{k+1}\|_1].$$

The updated sparse loading matrix  $\hat{A}_{SO}^{k+1}$  and specific variance matrix  $\hat{D}_{SO}^{k+1}$  are given by ( $i = 1, 2, \dots, p$ ):

$$\hat{A}_{SO}^{k+1} = (\sqrt{\lambda_1^{k+1}} v_{SO1}^{k+1}, \dots, \sqrt{\lambda_m^{k+1}} v_{SOm}^{k+1}) = (\hat{a}_{SOij}^{k+1})_{p \times m},$$

$$\hat{D}_{SO}^{k+1} = \text{diag}(\hat{\sigma}_1^2, \hat{\sigma}_2^2, \dots, \hat{\sigma}_p^2), \quad \hat{\sigma}_i^2 = s_{ii} - \sum_{j=1}^m \hat{a}_{soij}^{k+1^2}. \quad (2)$$

### C. SOPC for LFM

For the LFM in equation (1), the data matrix  $X$  is generated with Laplace distributed errors. Using the SOPC-derived sparse loading matrix, the updated parameters  $\hat{A}_{SO}^{k+1}$  and  $\hat{D}_{SO}^{k+1}$  follow equation (2) where

$$\hat{\sigma}_i^2 = X_{k+1}^T X_{k+1}, \quad i = 1, 2, \dots, p.$$

This completes the integration of SOPC with the LFM.

## IV. NUMERICAL ANALYSIS

### A. Preparation work

1) *Statistical index*: The efficacy of the SOPC algorithm is quantified through the Frobenius-norm-based Mean Squared Error (MSE) between the true and estimated factor loading matrices. In simulation studies, the true loading matrix  $A$  is predefined, enabling direct accuracy assessment. The MSE is defined as

$$MSE_{\hat{A}} = \frac{1}{p^2} \|A - \hat{A}\|_F^2$$

where  $\hat{A}$  denotes the estimated loading matrix.

2) *Parameter setting*: The simulated data  $X$  is generated according to the generative model in (1), with the following configurations:

- Parameter ranges: Location parameter  $\mu \in [0, 1000]$ , while factor loadings  $A$  are bounded within  $[-1, 1]$ ;
- Distribution specifications: Latent factors  $F \sim \mathcal{N}(0, I_{m \times m})$ , noise terms  $\epsilon \sim \text{Laplace}(0, 1)$ , and idiosyncratic variances  $D \in (0, 1)$ .

### B. Simulation 1: Baseline performance

We set  $n = 1000$ ,  $p = 10$ ,  $m = 5$  and generate the data matrix  $X$ . Table I summarizes the KMO and Bartlett's sphericity test results.

As evidenced in Table I, all KMO values exceed 0.60, and the sphericity hypothesis is rejected at  $p < 0.05$ . Both statistical criteria substantiate the appropriateness of LFM for dimensionality reduction.

TABLE I: KMO and Bartlett's test results

| Laplace Distribution       | KMO  | $\chi^2$ | df | P-value        | Suitable |
|----------------------------|------|----------|----|----------------|----------|
| Univariate Symmetric       | 0.69 | 1739.30  | 45 | 0              | YES      |
| Precision Parameterization | 0.62 | 1649.50  | 45 | $1.42e^{-316}$ | YES      |
| Univariate Asymmetric      | 0.70 | 2679.67  | 45 | 0              | YES      |

C. Simulation 2: Sensitivity analysis

1) Impact of sample size ( $n$ ): We set  $p = 10$ ,  $m = 5$ , and varying  $n \in \{2000, 3000, 4000, 5000, 6000\}$ .

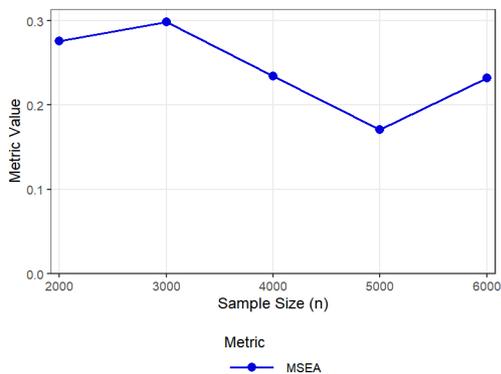


Fig. 1:  $MSE_{\hat{A}}$  values of SOPC for univariate symmetric model as  $n$  varies.

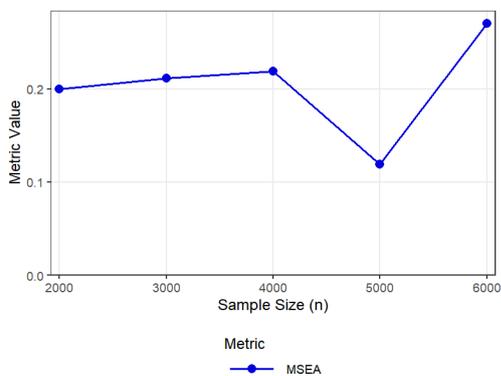


Fig. 2:  $MSE_{\hat{A}}$  values of SOPC for precision parameterization model as  $n$  varies.

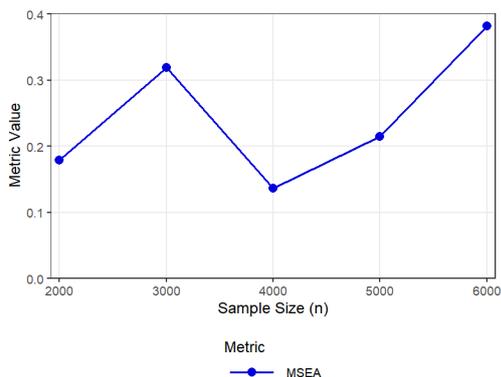


Fig. 3:  $MSE_{\hat{A}}$  values of SOPC for univariate asymmetric model as  $n$  varies.

- As shown in Fig. 1, the  $MSE_{\hat{A}}$  curve demonstrates strong stability, with values ranging from 0.17 to 0.29. The minimum is achieved at  $n = 5000$  ( $MSE_{\hat{A}} = 0.17$ ), followed by a slight increase at  $n = 6000$  ( $MSE_{\hat{A}} = 0.23$ ).

- Fig. 2 shows that  $MSE_{\hat{A}}$  values remain consistently stable within the range of 0.12 to 0.27, with the best performance observed at  $n = 5000$  ( $MSE_{\hat{A}} = 0.12$ ).
- In Fig. 3, moderate fluctuations are observed, with values ranging from 0.14 to 0.38. Notably, there is a significant increase in error at  $n = 6000$ , suggesting a sensitivity to larger sample sizes.

2) Impact of sample dimension ( $p$ ): We set  $n = 2000$ ,  $m = 5$ , and varying  $p \in \{10, 11, 12, 13, 14\}$ .

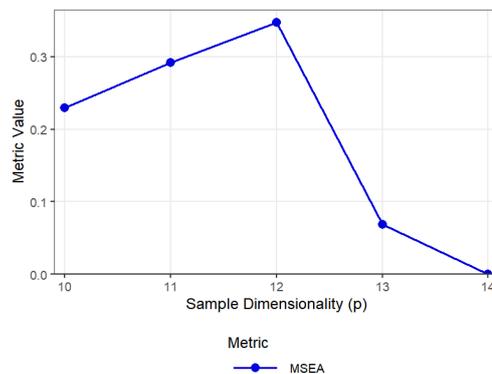


Fig. 4:  $MSE_{\hat{A}}$  values of SOPC for univariate symmetric model as  $p$  varies.

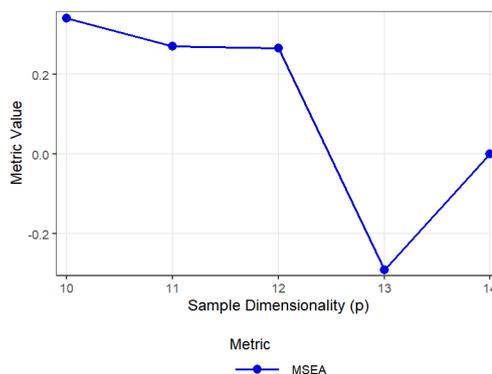


Fig. 5:  $MSE_{\hat{A}}$  values of SOPC for precision parameterization model as  $p$  varies.

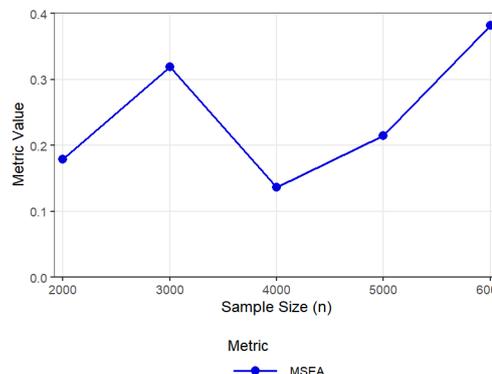


Fig. 6:  $MSE_{\hat{A}}$  values of SOPC for univariate asymmetric model as  $p$  varies.

- As shown in Fig. 4,  $MSE_{\hat{A}}$  peaks at  $p = 12$  ( $MSE_{\hat{A}} = 0.34$ ) before sharply declining to near-zero values at  $p = 14$ , suggesting a risk of overfitting with higher dimensions.
- Fig. 5 demonstrates that  $MSE_{\hat{A}}$  gradually decreases from 0.26 to 0.22, but exhibits negative values at  $p = 13$ , indicating instability in the estimation process at this point.
- In Fig. 6,  $MSE_{\hat{A}}$  fluctuates moderately within the range of 0.14 to 0.38, with the best performance observed at mid-range dimensions, specifically at  $p = 12$ .

D. Simulation 3: Comparative analysis

We compare the SOPC method against four alternatives: SPC, Projection Principal Component (PPC), Incremental Principal Component (IPC) and Stochastic Approximate Principal Component (SAPC). Parameters are fixed as  $p = 10$ ,  $m = 5$ ,  $\mu \in [0, 1000]$ ,  $A \in [-1, 1]$ ,  $F \sim \mathcal{N}_m(0, I_{m \times m})$ ,  $\varepsilon \sim \text{Laplace}(0, 1)$ , and  $D \in (0, 1)$ , with  $n \in \{2000, 3000, 4000, 5000, 6000\}$ .

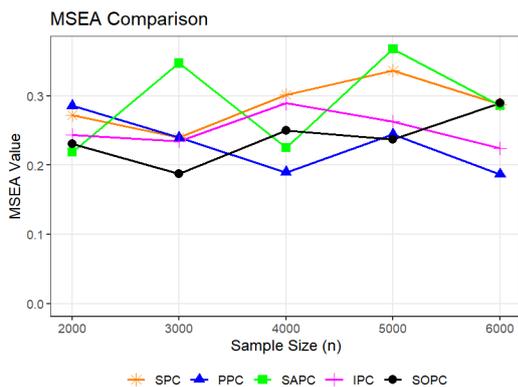


Fig. 7: Comparison of  $MSE_{\hat{A}}$  performance of SOPC, PPC, IPC, SPC, and SAPC methods for univariate symmetric model with the change of  $n$ -value.

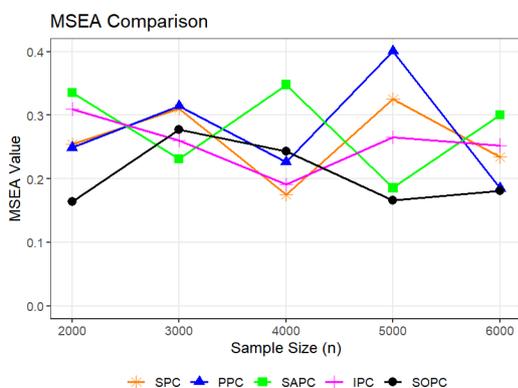


Fig. 8: Comparison of  $MSE_{\hat{A}}$  performance of SOPC, PPC, IPC, SPC, and SAPC methods for precision parameterization model with the change of  $n$ -value.

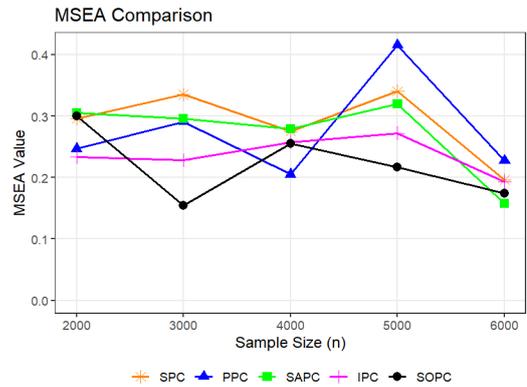


Fig. 9: Comparison of  $MSE_{\hat{A}}$  performance of SOPC, PPC, IPC, SPC, and SAPC methods for univariate asymmetric model with the change of  $n$ -value.

- As shown in Fig. 7, SOPC achieves the lowest  $MSE_{\hat{A}}$  values in the range of 0.19–0.24, outperforming SAPC (0.22–0.37) and all other methods.
- Fig. 8 demonstrates SOPC’s superior robustness, with its  $MSE_{\hat{A}}$  ranging from 0.16 to 0.28. In contrast, PPC shows considerable instability, with  $MSE_{\hat{A}}$  fluctuating between 0.19 and 0.40.
- In Fig. 9, SOPC consistently achieves the lowest error rates (0.15–0.20), while PPC exhibits the largest fluctuations, with  $MSE_{\hat{A}}$  spanning from 0.20 to 0.42.

These results underscore SOPC’s superior adaptability to variations in sample size compared to the other methods.

V. CONCLUSIONS AND SUGGESTIONS

The SOPC method exhibits high precision and stability in LFM, outperforming traditional principal component analysis. While sample size variations minimally affect performance, higher dimensions may increase estimation errors. Future work should integrate distributed computing for efficiency and extend the method to nonlinear models. This approach provides an effective solution for real-time analysis of complex datasets.

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