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# Joint Factors And Rank Canonical Estimation For The Polyadic Decomposition Based On Convex Optimization

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ABSTRACT Estimating the minimal number of rank-1 tensors in the Canonical Polyadic Decomposition (CPD) known as the canonical rank is a challenging area of research. To address this problem, we propose a method based on convex optimization to jointly estimate the CP factors and the canonical rank called FARAC for Joint FActors and RAnk canonical estimation for the PolyAdiC Decomposition. We formulate the FARAC method as a convex optimization problem in which a sparse promoting constraint is added to the superdiagonal of the core tensor of the CPD while the frobenius norm of the offdiagonal terms is constrained to be bounded. We propose an alternated minimization strategy of the Lagrangien to solve the proposed optimization problem. Compared to the state-of-the-art method, FARAC shows very good performance in terms of accuracy of the rank estimation for large range of values of SNRs. In addition, FARAC can handle the case where the canonical rank exceeds one of the dimensions of the input tensor. Several numerical experiments are conducted in the presence of various levels of noise.

INDEX TERMS Rank and factors estimation, Canonical Polyadic Decomposition, convex optimization

### I. INTRODUCTION

arge amounts of data are usually generated by sensor networks, massive experiment, simulations, *etc*. In a very wide range of applications, data need more than two dimensions to be efficiently described [2], [7], [21], [29]. To represent such data, multidimensionnal arrays (*a.k.a.* tensors) are suitable in order to capture complex interactions among input features of data. Tensors can be seen as a generalization of vectors (1st order tensors) and matrices (2nd order tensors). The order of a tensor is its number of dimensions. Various tensor decompositions exist to mitigate the curse of dimensionality *i.e.* to avoid the exponential growth of storage cost [3], [8], [14], [15], [22].

The most popular tensor decomposition is the Canonical Polyadic Decomposition (CPD) [13]. The CPD is widespreadly used in differents fields such as as chemometrics, telecommunications, blind source separation [11], [17]–[19], [27]. The CPD is particularly attractive due to its uniqueness property [1], [10]. However, the CPD requires the knowledge of the rank. Unfortunately, the problem of

determining the canonical rank is NP-hard [5].

In the present study, we propose estimating both of the canonical rank and the CP factors from noisy observations. We give a formulation of the problem of interest as a convex optimization problem. The reminder of the paper is organised as follows:

First, we introduce some notations and preliminaries in multilinear algebra in section II. We then review some existing works for the estimation of the canonical rank III. In section IV, we present our proposed approach and describe our algorithm for solving it. As a final step, we perform several numerical experiments in section VI to evaluate and compare our proposed approach with the CORCONDIA method.

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# II. NOTATIONS AND ALGEBRAIC BACKGROUND

In this section, we recall some algebraic definitions on tensor algebra from [13]:

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**Definition** 1: (Inner product): The inner product of two N-order tensors  $\mathcal{X}, \mathcal{Y} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$  is defined as:

$$\langle \mathcal{X}, \mathcal{Y} \rangle = \sum_{i_1=1}^{I_1} \sum_{i_2=1}^{I_2} \cdots \sum_{i_N=1}^{I_N} \mathcal{X}(i_1, \dots, i_N) \mathcal{Y}(i_1, \dots, i_N).$$

**Definition** 2: (Tensor Frobenius Norm): The norm of a tensor  $\mathcal{X}$  is defined as:

$$||\mathcal{X}||_F = \sqrt{\sum_{i_1=1}^{I_1} \sum_{i_2=1}^{I_2} \cdots \sum_{i_N=1}^{I_N} \mathcal{X}^2(i_1, \dots, i_N)}.$$

**Definition** 3: (n-mode multiplication) The n-mode product of a tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$  with a matrix  $U \in \mathbb{R}^{J \times I_n}$  is a tensor of order N and size  $I_1 \times \ldots I_{n-1} \times J \times \cdots \times I_N$  and is given for instance by:

$$(\mathcal{X} \times_n U)(i_1, \dots, i_{n-1}, i, i_{n+1}, \dots, i_N) = \sum_{i_n=1}^{I_n} \mathcal{X}(i_1, \dots, i_N)U(i, i_n).$$

**Definition** 4: (Diagonal tensor) A tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$  is diagonal if all of its entries are zero except the ones in its superdiagonal *i.e.*:  $\mathcal{X}(i_1,\ldots,i_N) \neq 0$  only if  $i_1=i_2=\cdots=i_N$ .

**Definition** 5: (Rank-one tensor) A tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$  of order N is rank one if it can be written as the outer product of N vectors:

$$\mathcal{X} = u_1 \circ \dots \circ u_N := \bigcap_{n=1}^N u_n, \tag{1}$$

where  $u_n \in \mathbb{R}^{I_n}$  with  $1 \le n \le N$ .

**Definition** 6: (Unfolding) The n-mode unfolding of a tensor  $\mathcal{X}$  is a matrix denoted by  $X_{(n)}$  where its columns are the n-mode fibers of  $\mathcal{X}$ .

**Definition** 7: (Tucker model) The Tucker decomposition transforms a tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$  of multilinear ranks  $(R_1, \ldots, R_N)$  using N factors  $U_n \in \mathbb{R}^{I_n \times R_n}$  by a core tensor  $\mathcal{G} \in \mathbb{R}^{R_1 \times \cdots \times R_N}$  as follows:

$$\mathcal{X} = \mathcal{G} \times_1 U_1 \cdots \times_N U_N := \mathcal{G} \times_{n=1}^N U_n$$
$$= \sum_{r_1}^{R_1} \cdots \sum_{i_N}^{R_N} \mathcal{G}(r_1, \dots, r_N) u_{1, r_1} \circ \cdots \circ u_{N, r_N},$$

where  $u_{n,r} \in \mathbb{R}^{I_n}$  is the r-th column of the n-th factor matrix  $U_n$  with  $1 \le n \le N$ .

**Definition** 8: (CANDECOMP Decomposition (CPD)) A tensor of order N and canonical rank R follows a CPD if it admits a factorization as a sum of R rank-one tensors. More formally, for a tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$ , the CPD is given by:

$$\mathcal{X} = \sum_{r=1}^{R} u_{1,r} \circ \dots \circ u_{N,r} = \sum_{r=1}^{R} \bigcap_{n=1}^{N} u_{n,r},$$
 (2)

CPD is a special case of the Tucker model where the core tensor is diagonal and the multilinear ranks are all equal to R. Hence, CPD in eq. (2) can be written in the Tucker format as follows:

$$\mathcal{X} = \mathcal{I} \times_{n=1}^{N} U_{n}.$$

**Definition** 9: (Tensor rank) The rank R of a tensor is defined as the smallest number of components in an exact CPD. For a 3-order tensor  $\mathcal{X} \in \mathbb{R}^{I_1 \times I_2 \times I_3}$ , we have the following upper bound of its maximum rank (the largest attainable

$$R \le \min\{I_1I_2, I_1I_3, I_2I_3\}.$$

### **III. RELATED WORKS**

rank) [10]:

Existing approaches for the canonical rank detection include:

- The CORe CONsistency DIAgnostic (CORCONDIA) [23] is a heuristic method to detect the number of components of the CP model. It measures the similarity between the estimated core tensor from the ideal identity core (called core consitency) for different CP models. By analyzing the gap between the core consistency of different CP models, it determines the correct number of components. Taking the last model (starting with the one-component model) whose core array is still similar to the ideal diagonal tensor gives the proper number of components to use. CORCONDIA is intuitive and a simple approach. However, its computation is prohibitive even for small tensors [20]. In addition, the CORCONDIA method can not handle the case when the canonical rank exceeds one of the dimensions of the input tensor.
- A fast version of CORCONDIA is presented in [20]. It proposes an efficient algorithm for computing the CORCONDIA diagnostic which exploits data sparsity and scales well as the tensor size increases. In cases where either the tensor or the factors or both are sparse [20], their algorithm significantly outperforms the state of the art baselines and scales well when the tensor size increases. In the fully dense scenario, their proposed algorithm is as good as the state of the art (The CORCONDIA method) for rank estimation.
- Automatic Relevance Determination (ARD) [16] is a
  bayesian approach applied for the Tucker and CP model.
  In ARD, entries of CP factors are assigned a Gaussian
  prior. The objective is to find the rank and the CP factors
  by solving an l<sub>2</sub>-regularized CP decomposition. By
  assigning priors to model hyperparameters and learning
  the hyperparameters of these priors, ARD turns excess
  of components and simplify the core structure.
- Detection of number of components in CANDECOMP models via minimum description length (N-D MDL) [12] is an extension of the minimum description length (MDL) to the multilinear case for detecting the number of components in a CP model using the generalized unfolding of the observation tensor. Under the condition that the tensor rank is smaller than the size of the most squared unfolded matrix, the generalized N-D MDL



criterion estimates the number of components of the CPD. The drawback of the multilinear MDL algorithm is that it fails to work when the tensor rank is larger than the size of the most squared unfolded matrix [12].

### IV. PROPOSED METHOD (FARAC)

The purpose of this section is to present the method FARAC for estimating the canonical rank and the CP factors simultaneously. This is achieved by minimizing the core tensors' superdiagonal using the  $l_0$  norm as an objective function, adding a constraint on the reconstruction error and an additional one on the offdiagonal terms by allowing them to be non-zero but bounded. Our goal is to find a CP core tensor structured as shown in Fig. 1. The canonical rank is the number of strictly positive and ordered values of the tensors' superdiagonal  $^1$ .

In the following,  $\mathcal{G}$  denotes the CP core tensor,  $\lambda$  its superdiagonal and  $\tilde{\mathcal{G}}$  is defined according to:

$$\mathcal{G} = \operatorname{diag}(\lambda) + \tilde{\mathcal{G}},$$

where  $diag(\lambda)$  is a diagonal tensor with  $\lambda$  in its superdiagonal. Our optimization problem can then be expressed mathematically as follows:

minimize 
$$\|\lambda\|_0$$
,  
subject to  $\frac{1}{2} \left\| \mathcal{X} - \mathcal{G} \underset{n=1}{\overset{N}{\times}_n} U_n \right\|_F^2 \le \epsilon_1$ , (3)  
 $\frac{1}{2} \left\| \tilde{\mathcal{G}} \right\|_F^2 \le \epsilon_2$ ,

where  $\epsilon_1$  and  $\epsilon_2$  are small positive constants. Unlike existing methods that estimate the canonical rank, we do not constrain the CP factors to be orthogonal [4], [9]. However, since minimizing the  $l_0$  norm is NP-hard [28], we minimize the  $l_1$  norm of  $\lambda$ . Eq. (3) becomes:

minimize 
$$\gamma_1 ||\lambda||_1$$
,

subject to  $\frac{1}{2} \left| \left| \mathcal{X} - \mathcal{G} \underset{n=1}{\overset{N}{\times}_n} U_n \right| \right|_F^2 \le \epsilon_1$ ,

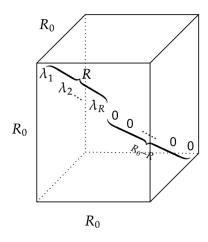
 $\frac{1}{2} \left| \left| \tilde{\mathcal{G}} \right| \right|_F^2 \le \epsilon_2$ ,

(4)

where  $\gamma_1$  is a strictly positive hyper-parameter.

### A. DERIVATION OF OUR APPROACH (FARAC)

In this section, we present the proposed approach for solving the optimization problem with constraints in eq. (4). Recall that the Lagrangien function is the augmented objective function by the constraint equations using the Lagrangien



**FIGURE 1.** CP core tensor  $\mathcal G$  of a rank-R tensor of order 3 and size  $R_0$  with  $\lambda_1>\cdots>\lambda_R>0$  .

multipliers. Following this, the Lagrangien function of the problem in eq. (4) is given by:

$$\mathcal{L}_{\mathcal{G},\{U_n\}} = \gamma_1 ||\lambda||_1 + \frac{\gamma_2}{2} \left( \left| \left| \mathcal{X} - \mathcal{G} \underset{n=1}{\overset{N}{\times}_n} U_n \right| \right|_F^2 - \epsilon_1 \right) + \frac{\gamma_3}{2} \left( \left| \left| \tilde{\mathcal{G}} \right| \right|_F^2 - \epsilon_2 \right),$$
(5)

where  $\gamma_2$  and  $\gamma_3$  are two strictly positive Lagrange multipliers. Following the Lagrangien method [24],  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  is minimized w.r.t  $\{U_n\}_n$ ,  $\tilde{\mathcal{G}}$  and  $\lambda$ . To do that, we will proceed iteratively. At each iteration, we first minimize  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t the n-th CP factor  $U_n$  assuming the remaining factors and  ${\cal G}$  are known. This is a classical linear regression problem. Next, we minimize  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\tilde{\mathcal{G}}$  using the factors  $\{U_n\}_n$ from the previous step. The convexity w.r.t  $\tilde{\mathcal{G}}$  is demonstrated in Appendix B. Then, we update  $\tilde{\mathcal{G}}$  using the Adam optimizer [6]. As a final step, we minimize  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\lambda$ , which is also a convex optimization problem as shown in Appendix B. Its solution is given by the soft thresholding operator [26] using the current updates of  $\tilde{\mathcal{G}}$  and  $\{U_n\}_n$ . We provide in the following, the final expressions of the exact solutions of CP factors, the gradient of the Lagrangien w.r.t the  $\hat{\mathcal{G}}$  and the formula to update the superdiagonal of  $\mathcal{G}$ . Details of the computations can be found in Appendix A.

• CP factors  $U_n$ :

$$U_n = X_{(n)} \left( \bigotimes_{\substack{l=N \\ l \neq n}}^{l=1} U_l \right) G_{(n)}^T \left( \bigotimes_{\substack{l=N \\ l \neq n}}^{l=1} \left( U_l^T U_l \right)^{-1} \right). \quad (6)$$

ullet Gradient of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t the offdiagonals of  $\mathcal{G}$  i.e

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<sup>&</sup>lt;sup>1</sup>The sign of the superdiagonal entries can always be absorbed into the factors.

 $\mathcal{G}(r_1,\ldots,r_N)$  such that  $r_1\neq r_2$  or  $\ldots$  or  $r_{N-1}\neq r_N$ :

$$[\nabla_{\tilde{\mathcal{G}}}(\mathcal{L}_{\mathcal{G},\{U_n\}})](r_1,\ldots,r_N) =$$

$$-\gamma_2 \sum_{i_1=1}^{I_1} \cdots \sum_{i_N}^{I_N} \mathcal{A}(i_1,\ldots,i_N)$$

$$\left[\prod_{n=1}^N U_n(i_n,r_n)\right] + \gamma_3 \tilde{\mathcal{G}}(r_1,\ldots,r_N), \quad (7)$$

where 
$$\mathcal{A}=\mathcal{X}-\mathcal{G}\overset{N}{\underset{n=1}{\times}_{n}}U_{n}.$$
   
 • Formula to update the superdiagonal of  $\mathcal{G}$ :

$$\lambda_{l} = S_{\frac{\gamma_{1}}{\gamma_{2}}} \left( -\left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{\substack{r=1\\r\neq l}}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle + \left\langle \mathcal{X} - \tilde{\mathcal{G}} \times_{n}^{N} U_{n}, \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle \right),$$
(8)

where  $S_{\frac{\gamma_1}{2}}$  is the soft thresholding operator [26] defined

$$S_{\mu}(x) = \begin{cases} x - \mu & \text{if } x > \mu \\ 0 & \text{if } |x| \le \mu \\ x + \mu & \text{if } x < -\mu \end{cases}$$
 (9)

The whole algorithm of our derived approach is described in Algorithm 1.

## V. COMPLEXITY ANALYSIS

• We evaluate the complexity analysis of the Algorithm 1 by taking into consideration the svds in the initialization and the main parts of the algorithm as the computations of the different gradients and the shrinkage operator. The complexity is evaluated as follows:

$$\mathcal{O}\left(\left(\sum_{n=1}^{N} I_{n}\right) R_{0}^{2} + T \left[N + \left(\sum_{n=1}^{N} I_{n}\right) R_{0} + \left(\prod_{n=1}^{N} I_{n}\right) \left(N + R_{0} + N R_{0}^{N-1}\right)\right]\right), \quad (10)$$

where T is the number of iterations.

• If  $I = \max(I_1, \dots, I_N)$ , the complexity in eq. (10)

$$\mathcal{O}\left(IR_0^2 + T\left(N + IR_0 + I(N + R_0 + NR_0^{N-1})\right)\right).$$

### VI. NUMERICAL EXPERIMENTS

### A. PROTOCOL EXPERIMENT

Using the CP model, we create a synthetic rank-R real valued tensor  $\mathcal{X}$  of order N. The CP factors are derived from a single normal standard distribution.  $\mathcal{B}$  is a zero mean, unit variance and white noise added so that we use different values of SNR (Signal to Noise Ratio) ranging from 0db to 40db. Our noisy data are then obtained by:

$$\mathcal{X}_{\text{noise}} = \mathcal{X}' + \sigma \mathcal{B}', \tag{11}$$

Algorithm 1: Joint FActors and RAnk Canonical estimation for the Polyadic decomposition (FARAC)

**Input:**  $\mathcal{X} \in \mathbb{R}^{I_1 \times ... I_N}$ : CP Tensor,

 $R_0$ : Upper bound of the rank of  $\mathcal{X}$ .

**Require:**  $\alpha$ : Stepsize;

 $\beta_1, \beta_2 \in [0, 1]$ : Exponential decay rates;

T: Maximum number of iterations

 $\epsilon = 10^{-8}$ : Parameter to avoid numerical instability.

**Initialize:** For  $1 \le n \le N$ :

$$U_n^{(0)} = \left\{ \begin{array}{l} svd(X_{(n)},R_0) \text{ if } I_n > R_0, \\ \operatorname{conc} \left( svd(X_{(n)},R_0), \\ (R_0 - I_n) \text{random uniform vectors} \right) \text{ else.} \end{array} \right.$$

- $\mathcal{G}^{(0)} \sim \mathcal{U}(0,1)$  of order N and size  $(R_0,\ldots,R_0)$ .  $m_{\tilde{\mathcal{G}}}^{(0)} = v_{\tilde{\mathcal{G}}}^{(0)} = 0$  (Initialize the first and the second moment estimates ).

**for** t = 1, ..., T:

- 1: Compute  $U_n^{(t)}$  from eq. (6). 2: Compute the gradients of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\tilde{\mathcal{G}}$  using (7).
- 3: Update biased first moment estimate of the offdiags:

$$m_{\tilde{q}}^{(t)} \leftarrow \beta_1 m_{\tilde{q}}^{(t-1)} + (1 - \beta_1) \nabla_{\tilde{q}}^{(t)}$$

4: Update biased second raw moment estimate of the offdiags:

$$v_{\tilde{\mathbf{G}}}^{(t)} \leftarrow \beta_2 v_{\tilde{\mathbf{G}}}^{(t-1)} + (1 - \beta_2) \nabla^{2}_{\tilde{\mathbf{G}}}^{(t)},$$

5: Compute bias-corrected first and second moment estimates of the offdiags:

$$\hat{m}_{\tilde{\mathcal{G}}}^{(t)} \leftarrow \frac{m_{\tilde{\mathcal{G}}}^{(t)}}{1 - \beta_1^t}; \quad \hat{v}_{\tilde{\mathcal{G}}}^{(t)} \leftarrow \frac{v_{\tilde{\mathcal{G}}}^{(t)}}{1 - \beta_2^t},$$

6: Update  $\hat{\mathcal{G}}$ :

$$\tilde{\mathcal{G}}^{(t)} \leftarrow \tilde{\mathcal{G}}^{(t-1)} - \alpha \frac{\hat{m}_{\tilde{\mathcal{G}}}^{(t)}}{\sqrt{\hat{v}_{\tilde{\mathcal{G}}}^{(t)}} + \epsilon}.$$

7: Update the superdiagonal of  $\mathcal{G}$  using eq. (8):

$$\begin{split} \lambda_l^{(t)} &= S_{\frac{\gamma_1}{\gamma_2}} \Bigg( - \left\langle \bigcap_{n=1}^N u_n^{(t)}, \sum_{\substack{r=1\\r \neq l}}^{R_0} \lambda_r^{(t)} \bigcap_{n=1}^N u_n^{(t)} \right\rangle + \\ & \left\langle \mathcal{X} - \tilde{\mathcal{G}} \mathop{\times}_{n=1}^N U_n^{(t)}, \bigcap_{n=1}^N u_n^{(t)} \right\rangle \Bigg), \end{split}$$

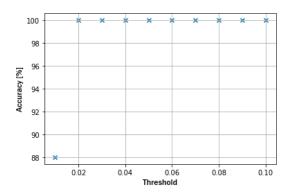
where  $S_{\frac{\gamma_1}{\gamma_2}}$  is defined in eq. (9).

end for

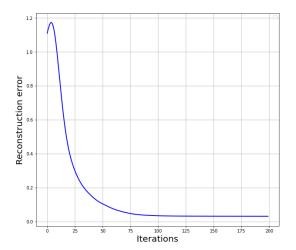
CP-rank: Number of non-zero values of the superdiagonal of  $\mathcal{G}$ .

**CP factors**: Columns of  $U_n^{(T)}$  with indices of non-zero values of the superdiagonal of  $\mathcal{G}$ 

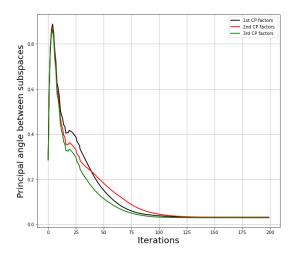
**Returns:** [CP-rank, CP factors]



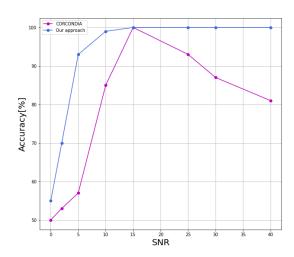
**FIGURE 2.** Accuracy of rank estimation of a tensor of size  $5\times5\times5$  w.r.t the threshhold parameter. The true rank is R=2 and  $R_0=5$ .



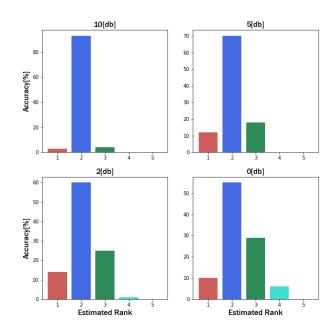
**FIGURE 3**. This figure shows the convergence curve of the mean reconstruction error using the Relative Square Error (RSE) along iterations using FARAC. We have used a noisy tensor  $\mathcal{X}_{\text{noise}}$  of size  $5\times5\times5$  with SNR = 25db. The threshold paramter is equal to 0.02.



**FIGURE 4.** This figure shows the mean principal angle between the three subspaces spanned by CP factors along iterations using our method. We have used a noisy tensor  $\mathcal{X}_{\mathsf{noise}}$  of size  $5 \times 5 \times 5$  with SNR = 25db. The threshold parameter is equal to 0.02.



**FIGURE 5.** Accuracy of FARAC Vs. CORCONDIA w.r.t SNR for a tensor of a noisy tensor  $\mathcal{X}_{\text{noise}}$  with SNR values ranging from 0db to 40db.  $\mathcal{X}_{\text{noise}}$  is of size  $5\times5\times5$ . Different CP models with ranks ranging from 1 to 5 are used to fit CORCONDIA. The true rank is R=2.

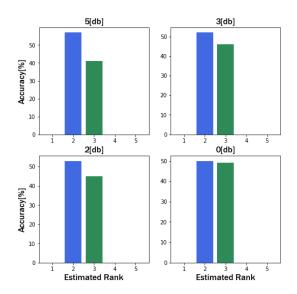


**FIGURE 6.** Accuracy of rank estimation using FARAC of a noisy tensor  $\mathcal{X}_{\mathsf{noise}}$  with different low SNR values ranging from 0db to 10db.  $\mathcal{X}_{\mathsf{noise}}$  is of size  $5 \times 5 \times 5$ . The large bound of rank used is  $R_0 = 5$  and the true rank is R = 2.

SNR [db]	15	20	25	30	35	40
Accuracy [%]	72	77	87	91	96	100

**TABLE 1.** Accuracy of FARAC w.r.t SNR for a tensor of a noisy tensor  $\mathcal{X}_{\text{noise}}$  of size  $5\times5\times5$ . The true rank is R=6 and  $R_0=7$ .

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**FIGURE 7.** Accuracy of the CORCONDIA approach for a noisy tensor  $\mathcal{X}_{\mathsf{noise}}$  with SNR values ranging from 0db to 5db.  $\mathcal{X}_{\mathsf{noise}}$  is of size  $5 \times 5 \times 5$ . Different CP models with ranks ranging from 1 to 5 are used to fit CORCONDIA. The true rank is R=2

where  $\mathcal{X}' = \frac{\mathcal{X}}{||\mathcal{X}||_F}$  and  $\mathcal{B}' = \frac{\mathcal{B}}{||\mathcal{B}||_F}$ . Hence, the SNR will be calculated using the following formula:

$$SNR = -10\log_{10}\sigma^2.$$

FARAC has been validated on a rank-2 tensor  $\mathcal{X}$  of size  $I \times I \times I$  where  $R_0 = I = 5$ . Experiments will be conducted on a tensor with these parameters until other settings are indicated. For tensors with other orders and sizes, similar results are obtained. By using  $\mathcal{X}$  and eq. (11), we generate 100 noisy realisation of the inpout tensor. The accuracy is defined as the proportion of realisations where the estimated rank is accurate.

### B. RESULTS

- Fig. 2 illustrates the accuracy of estimated ranks using the proposed approach. We can clearly see that
  the FARAC approach can estimate the exact canonical
  rank while being robust to the choise of the threshold
  parameter.
- Fig. 3 shows the convergence curves of our method in the sense of the Reconstruction Square Error (RSE) at each iteration t given by:

$$\frac{||\mathcal{X}_{\text{noise}} - \mathcal{X}^{(t)}||_F}{||\mathcal{X}_{\text{noise}}||_F}.$$

We can see that the reconstruction error quickly decreases to zero w.r.t iterations.

• In Fig. 4, the recovery of the CP factors can be shown by checking the subspace error. This can be done by computing the principal angle [25] between the true subspaces and the estimated ones at each iteration t:

$$\theta^{(t)} = \arcsin\left(\sqrt{2}||U_nU_n^{\dagger} - \hat{U}_n^{(t)}\hat{U}_n^{(t)\dagger}||_F\right),\,$$

- where  $U_n$  and  $\hat{U}_n^{(t)}$  are respectively the exact and the estimated factors at iteration t. We see that the principal angle between the subspaces of the three CP factors converges to a low value w.r.t iterations.
- In Fig. 5, we compare the FARAC method to the COR-CONDIA method w.r.t SNR values. We can see that FARAC clearly outperforms the state-of-the-art method for large range values of SNRs.
- In Table 1, we show the accuracy of rank estimation using FARAC when the true rank exceeds one of the dimensions of the input tensor. As we can see, FARAC can handle this difficult scenario for a large range of SNR values. On the other hand, the CORCONDIA approach cannot deal with that case [12].
- According to Fig. 6, we can say that at very low SNR, FARAC tends to overestimate the true rank. In contrast, rank estimation using the CORCONDIA method becomes a very difficult task at low SNR according to Fig. 7.

Compared with the state-of-the-art method, FARAC shows very good performance in terms of the accuracy of rank estimation for large range of values of SNRs while being robust to the choise of the threshold parameter. Moreover, FARAC deals with the difficult case where the rank exceeds one of the dimensions of the input tensor.

### VII. CONCLUSION

We have addressed in this work the challenging problem of the canonical rank estimation by defining it as a convex optimization problem. The proposed method called FARAC jointly estimates the canonical rank and the CP factors. We have compared FARAC to the well-known CORCONDIA method and found that FARAC is much more accurate. We have also shown that FARAC shows strong robustness to the choise of the threshold parameter and can handle the difficult case when the rank exceeds one of the dimensions of the tensor unlike to the CORCONDIA method. Applications such as image denoising could be explored using the FARAC method.

# A. DETAIL COMPUTATIONS FOR THE DERIVATION OF FARAC

Let us recall the Lagrangien of the problem (4) that we want to minimize w.r.t  $(U_n)_n$ ,  $\tilde{\mathcal{G}}$  and  $\lambda$ :

$$\mathcal{L}_{\mathcal{G},\{U_n\}} = \gamma_1 ||\lambda||_1 + \frac{\gamma_2}{2} \left( \left| \left| \mathcal{X} - \mathcal{G} \underset{n=1}{\overset{N}{\times}_n} U_n \right| \right|_F^2 - \epsilon_1 \right) + \frac{\gamma_3}{2} \left( \left| \left| \tilde{\mathcal{G}} \right| \right|_F^2 - \epsilon_2 \right),$$

• We denote by  $\mathcal{L}_{U_n}$ , the part of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  that only depends on  $U_n$ .

$$\mathcal{L}_{U_n} = \frac{\gamma_2}{2} \left\| \mathcal{X} - \mathcal{G} \underset{n=1}{\overset{N}{\times}} U_n \right\|_F^2$$
 (12)



The matricized form of eq. (12) is given by:

$$\mathcal{L}_{U_n} = \left| \left| X_{(n)} - U_n G_{(n)} \left( \bigotimes_{\substack{l=N \\ l \neq n}}^{l=1} U_l^T \right) \right| \right|_F^2,$$

As a difference from the derivation of a classical CPD,  $G_{(n)}$  is not the identity matrix.

The gradient of  $\mathcal{L}_{U_n}$  w.r.t  $U_n$  can be derived as in a classical matrix linear regression problem:

$$\nabla_{U_n}(\mathcal{L}_{\mathcal{G},\{U_n\}}) = \gamma_2 \left( -X_{(n)} \left( \bigotimes_{\substack{l=N\\l \neq n}}^{l=1} U_l \right) G_{(n)}^T + U_n \left( \bigotimes_{\substack{l=N\\l \neq n}}^{l=1} U_l^T U_l \right) \right)$$

We set  $\nabla_{U_n}(\mathcal{L}_{\mathcal{G},\{U_n\}})$  to 0 and get the exact formula of  $U_n$ :

$$U_n = X_{(n)} \left( \bigotimes_{\substack{l=N\\l \neq n}}^{l=1} U_l \right) G_{(n)}^T \left( \bigotimes_{\substack{l=N\\l \neq n}}^{l=1} \left( U_l^T U_l \right)^{-1} \right).$$

We denote by \(\mathcal{L}\_{\tilde{\mathcal{G}}}\), the part of \(\mathcal{L}\_{\mathcal{G},\{U\_n\}}\) that only depends on \(\tilde{\mathcal{G}}\):

$$\mathcal{L}_{\tilde{\mathcal{G}}} = \frac{\gamma_2}{2} \sum_{i_1, \dots, i_N} \underbrace{\left(\mathcal{X}(i_1, \dots, i_N) - \left(\mathcal{G} \underset{n=1}{\overset{N}{\times}} U_n\right) (i_1, \dots, i_N)\right)^2}_{\mathcal{A}^2(i_1, \dots, i_N)} + \frac{\gamma_3}{2} \sum_{r_1, \dots, r_N} \underbrace{\tilde{\mathcal{G}}^2(r_1, \dots, r_N)}_{\mathcal{G}^2(r_1, \dots, r_N)}$$

Let us derive  $\mathcal{L}_{\tilde{\mathcal{G}}}$  w.r.t  $\tilde{\mathcal{G}}(r_1,\ldots,r_N)$  such that  $r_1 \neq r_2$  or  $\ldots$  or  $r_{N-1} \neq r_N$  (diagonal elements are excluded since they are equal to 0):

$$[\nabla_{\tilde{\mathcal{G}}} \mathcal{L}_{\tilde{\mathcal{G}}}](r_1, \dots, r_N) = -\gamma_2 \sum_{i_1, \dots, i_N} \mathcal{A}(i_1, \dots, i_N) \left( \prod_{n=1}^N U_n(i_n, r_n) \right) + \gamma_3 \tilde{\mathcal{G}}(r_1, \dots, r_N).$$
(13)

• We want to derive  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\lambda$ . Let us first rewrite it as follows:

$$\mathcal{L}_{\mathcal{G},\{U_n\}} = \gamma_1 ||\lambda||_1 + \frac{\gamma_2}{2} \left\| \mathcal{X} - \left( \tilde{\mathcal{G}} + diag(\lambda) \right) \sum_{n=1}^N U_n \right\|_F^2 + \frac{\gamma_3}{2} \left\| \tilde{\mathcal{G}} \right\|_F^2,$$

$$= \gamma_1 ||\lambda||_1 + \frac{\gamma_2}{2} \left\| \left( \mathcal{X} - \tilde{\mathcal{G}} \sum_{n=1}^N U_n \right) - \sum_{n=1}^{R_0} \lambda_r \bigcap_{n=1}^N u_n^{(r)} \right\|_F^2 + \frac{\gamma_3}{2} \left\| \tilde{\mathcal{G}} \right\|_F^2$$

 We denote by L<sub>λ</sub>, the part of L<sub>G,{U<sub>n</sub>}</sub> that only depends on λ. L<sub>λ</sub> is given as follows:

$$\mathcal{L}_{\lambda} = \gamma_{1} ||\lambda||_{1} + \frac{\gamma_{2}}{2} \left( \left\| \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\|_{F}^{2} - 2 \left\langle \mathcal{X} - \tilde{\mathcal{G}} \bigcap_{n=1}^{N} U_{n}, \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle + \left\| \mathcal{X} - \tilde{\mathcal{G}} \bigcap_{n=1}^{N} U_{n} \right\|_{F}^{2} \right)$$

$$= \gamma_{1} \left( |\lambda_{l}| + \sum_{r \neq l} |\lambda_{r}| \right) + \frac{\gamma_{2}}{2} \left( \left\| \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\|_{F}^{2} + \lambda_{l}^{2} \left\| \bigcap_{n=1}^{N} u_{n}^{(l)} \right\|_{F}^{2} + 2 \left\langle \lambda_{l} \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle - 2 \left\langle \mathcal{X} - \tilde{\mathcal{G}} \bigcap_{n=1}^{N} U_{n}, \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle + \left\| \mathcal{X} - \tilde{\mathcal{G}} \bigcap_{n=1}^{N} U_{n}, \lambda_{l} \bigcap_{n=1}^{N} U_{n} \right\|_{F}^{2} \right). \quad (14)$$

Let us derive  $\mathcal{L}_{\lambda}$  w.r.t  $\lambda_l$  and set it to 0.

$$\nabla_{\lambda_{l}} (\mathcal{L}_{\lambda}) = 0$$

$$\Leftrightarrow \gamma_{1} \frac{\partial |\lambda_{l}|}{\lambda_{l}} + \gamma_{2} \left( \lambda_{l} \left\| \bigcap_{n=1}^{N} u_{n}^{(l)} \right\|_{F}^{2} + \left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle$$

$$- \left\langle \mathcal{X} - \tilde{\mathcal{G}} \times_{n}^{N} U_{n}, \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle = 0$$

$$\Leftrightarrow \lambda_{l} + \frac{\gamma_{1}}{\gamma_{2} \left\| \bigcap_{n=1}^{N} u_{n}^{(l)} \right\|_{2}^{2}} \frac{\partial |\lambda_{l}|}{\partial \lambda_{l}} =$$

$$- \gamma_{2} \left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{r=1}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle$$

$$+ \gamma_{2} \left\langle \mathcal{X} - \tilde{\mathcal{G}} \times_{n}^{N} U_{n}, \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle. \quad (15)$$

Furthermore, we have the following:

$$\left| \left| \bigcap_{n=1}^{N} u_n^{(l)} \right| \right|^2 = \prod_{n=1}^{N} \left| \left| u_n^{(l)} \right| \right|^2.$$

Since  $u_n^{(l)}$  are unit vectors, we have  $\left\| \bigcap_{n=1}^N u_n^{(l)} \right\|^2 = 1$ . Hence we find the following:

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$$\lambda_{l} + \frac{\gamma_{1}}{\gamma_{2}} \frac{\partial |\lambda_{l}|}{\partial \lambda_{l}} = -\left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{\substack{r=1\\r\neq l}}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle + \left\langle \mathcal{X} - \tilde{\mathcal{G}} \times_{n}^{N} U_{n}, \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle$$

Hence, using the soft thresholding operator  $S_{\frac{\gamma_1}{\gamma_2}}$ , we conclude the value of  $\lambda_l$ :

$$\lambda_{l} = S_{\frac{\gamma_{1}}{\gamma_{2}}} \left( -\left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{\substack{r=1\\r\neq l}}^{R_{0}} \lambda_{r} \bigcap_{n=1}^{N} u_{n}^{(r)} \right\rangle + \left\langle \mathcal{X} - \tilde{\mathcal{G}} \times_{n}^{N} U_{n}, \bigcap_{n=1}^{N} u_{n}^{(l)} \right\rangle \right)$$

### B. CONVEXITY

In this section, we will demonstrate that the Lagrangien in eq. (14) is convex w.r.t to  $\lambda$  and  $\tilde{\mathcal{G}}$  so that the global minimum will be reached. To do that, we will show that the hessian w.r.t  $\lambda$  and  $\tilde{\mathcal{G}}$  are positive semi-definite matrix.

• Convexity w.r.t  $\lambda$ : Let  $H_{\lambda}$  be the hessian of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\lambda$ . Using the gradient computed in eq. (15), we have:

$$H_{\lambda}(s,l) := \frac{\partial^2 \mathcal{L}_{\mathcal{G},\{U_n\}}}{\partial \lambda_s \partial \lambda_l} = \gamma_2 \left\langle \bigcap_{n=1}^N u_n^{(l)}, \bigcap_{n=1}^N u_n^{(s)} \right\rangle.$$

Let  $x \in \mathbb{R}^{R_0}$ ,

$$x^{T} H_{\lambda} x = \gamma_{2} \sum_{l,s} x_{l} x_{s} \left\langle \bigcap_{n=1}^{N} u_{n}^{(l)}, \bigcap_{n=1}^{N} u_{n}^{(s)} \right\rangle$$
$$= \gamma_{2} \left\langle \sum_{l} x_{l} \bigcap_{n=1}^{N} u_{n}^{(l)}, \sum_{s} x_{s} \bigcap_{n=1}^{N} u_{n}^{(s)} \right\rangle$$
$$= \gamma_{2} \left\| \sum_{l} x_{l} \bigcap_{n=1}^{N} u_{n}^{(l)} \right\|^{2} \geq 0,$$

since  $\gamma_2$  is a positive Lagrange multiplier.

• Convexity w.r.t  $\tilde{\mathcal{G}}$ :

We first put the elements  $\tilde{\mathcal{G}}(r_1,\ldots,r_N)$  in a vector  $x\in\mathbb{R}^{R_0^N-R_0}$  and use the same method as for  $\lambda$ .

Let  $H_{\tilde{\mathcal{G}}}$  be the hessian of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$  w.r.t  $\tilde{\mathcal{G}}$ . Using (13), we compute the second derivates of  $\mathcal{L}_{\mathcal{G},\{U_n\}}$ . The following derivatives are done only on the offdiagonals of  $\tilde{\mathcal{G}}$  since its diagonal is zero.

$$H_{\tilde{\mathcal{G}}}(r_1 \dots r_N, r'_1 \dots r'_N) := \frac{\partial^2 \mathcal{L}_{\mathcal{G}, \{U_n\}}}{\partial \tilde{\mathcal{G}}(r_1, \dots, r_N) \partial \tilde{\mathcal{G}}(r'_1, \dots, r'_N)}$$

$$= \gamma_2 \sum_{i_1, \dots, i_N} \left[ \prod_{n=1}^N U_n(i_n, r'_n) \right] \left[ \prod_{n=1}^N U_n(i_n, r_n) \right] + \gamma_3.$$

$$\begin{split} & \operatorname{Let} x \in \mathbb{R}^{R_0^N - R_0}, \\ & x^T H_{\tilde{\mathcal{G}}} x = \\ & \gamma_2 \sum_{i_1, \dots, i_N} \underbrace{\sum_{r_n} \left[ \prod_{n=1}^N U_n(i_n, r_n) \right]}_{\mathcal{Z}(i_1, \dots, i_N)} \underbrace{\sum_{r'_n} \left[ \prod_{n=1}^N U_n(i_n, r'_n) \right]}_{\mathcal{Z}(i_1, \dots, i_N)} \\ & + \gamma_3 x^T x \\ & = \gamma_2 ||\mathcal{Z}||^2 + \gamma_3 ||x||^2 \geq 0, \end{split}$$

since  $\gamma_2$  and  $\gamma_3$  are strictly positive values.

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