Journal homepage: http://iieta.org/journals/isi

A Novel Log-Based Tensor Completion Algorithm

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https://doi.org/10.18280/isi.250202	ABSTRACT
Received: 17 December 2019	In recent years, tensor completion problem, as a higher order generalization of matrix

Accepted: 17 December 2019 Accepted: 29 January 2020

Keywords:

tensor completion, log function of tensor, image inpainting, tensor decomposition

In recent years, tensor completion problem, as a higher order generalization of matrix completion, has received much attention from scholars engaged in computer vision, data mining, and neuroscience. The problem is often solved by convex relaxation, which minimizes the tensor nuclear norm instead of the n-rank of the tensor. However, tensor nuclear minimizes all the singular value at the same level, which is unfair to large singular values. To solve the problem, this paper defines a log function of tensor, and uses it as an approximation of tensor rank function. Then, a simple yet efficient log-based algorithm for tensor completion (Log-TC) was proposed to recover an underlying low n-rank tensor. The Log-TC was verified through experiments on randomly generated tensors and color image inpainting, in comparison with two tensor completion algorithms: fixed point iterative method for low-rank tensor completion (FP-LRTC) and fast low rank tensor completion algorithm (FaLRTC). The results show that our algorithm greatly outperformed the two contrastive methods.

1. INTRODUCTION

In the era of data explosion, tensors gain more and more attention thanks to their excellence in information preservation and structural properties. Tensors can be viewed as natural higher-order generalization of vectors and matrices. Therefore, many questions of tensors can be generalized from vectors and matrices.

Tensor completion, a natural generalization of matrix completion, has long been a research hotspot. As an imputation method for missing data, tensor completion aims to estimate missing data from very limited information of observed data. To date, this problem has been widely applied in recommendation system [1-7], signal processing [5-8], computer vision [9-13], and multi-relation link prediction [14-16].

Tensor decomposition is an important step of tensor completion. The main forms of tensor decomposition include CANDECOMP/PARAFAC (CP) decomposition [17, 18], Tucker decomposition [19-21], and Tensor-Train decomposition [22-24]. Our results are mainly based on Tucker decomposition.

In recent years, lots of algorithms have been developed to complete high-order tensors. Most of them adopts the lowrank structure assumption. It is known to all that tensor rank minimization is a non-deterministic polynomial-time (NP) hard problem. Therefore, many norms are defined as the convex surrogates of tensor rank. For example, Liu et al. [10] defined the trace norm for tensors, and proposed a completion strategy for low-rank tensors, laying the theoretical basis for low n-rank tensor completion.

Since all the singular values are simultaneously minimized, the convex surrogates may be insufficient to approximate the rank function. Therefore, many scholars have attempted to prove the advantage of approximating the rank for matrices, using nonconvex surrogate functions [25-28].

Recently, new results have been achieved in tensor completion through the extension of matrix situation. For instance, Huang et al. [29] put forward a tensor n-mode matrix unfolding truncated nuclear norm minimization, and solved it by the alternating direction method of multipliers. Han et al. [30] also presented an algorithm based on tensor truncated nuclear norm, and added a sparse regularization term into the objective function. Xue et al. [31] also carried out similar research.

In general, large singular values represent the significant information of interest, which should not be penalized excessively. By contrast, small singular values represent the noisy information, which should be penalized as zeros. All the singular values are minimized to the same level by tensor nuclear norm. This is obviously unfair to large singular values, which contain much more important information than small singular values. To solve the problem, log function was introduced to strike a balance between rank minimization and nuclear norm.

Taking the n-rank of a tensor as a sparsity measure, this paper solves the low-n-rank tensor completion problem in the following manner. A log function of tensor was defined, and used as an approximation of tensor rank function. In this way, the objective function in our optimization model become a nonconvex function, which is difficult to solve. Thus, the DC programming was adopted to solve the model. Next, a simple yet efficient log-based algorithm for tensor completion (Log-TC) was designed to recover an underlying low n-rank tensor. The performance of Log-TC algorithm was verified by computational results on recovery of synthetic data and color image.

The remainder of this paper is organized as follows: Section 2 briefly introduces tensor completion; Section 3 proposes an iterative scheme to minimize log function, and presents the



Log-TC algorithm; Section 4 compares our algorithm on randomly generated tensors and color image inpainting, with two low-rank tensor completion algorithms; Section 5 puts forward the conclusions.

2. PRELIMINARIES

2.1 Notations and definitions

Throughout this paper, each scalar is denoted by a normal letter, e.g. a, b, c, ...; each vector is denoted by a boldfaced lower-case letter, e.g. **a**, **b**, **c**, ...; each matrix is denoted by an upper-case letter, e.g. A, B, C, ...; each tensor is written as an italic letter, e.g. an N-order tensor is denoted as $\mathcal{X} \in \mathbb{R}^{l_1 \times \cdots \times l_N}$, where I_k is the dimensional size corresponding to mode-k, $k=1, \ldots, N$; each element of tensor \mathcal{X} is denoted as $x_{i_1 \cdots i_N} \in \mathcal{X}$.

The Frobenius norm of \mathcal{X} is defined as the square root of the inner product of twofold tensors:

$$\|\mathcal{X}\|_F = \sqrt{\langle \mathcal{X}, \mathcal{X} \rangle} = \sqrt{\sum_{i_1=1}^{I_1} \cdots \sum_{i_N=1}^{I_N} x_{i_1 \cdots i_N}^2}$$
(1)

The CP decomposition factorizes the target tensor into a sum of component one-rank tensors. For example, a third-order tensor $\mathcal{X} \in \mathbb{R}^{I \times J \times K}$ can be written as:

$$\mathcal{X} \approx \sum_{r=1}^{R} a_r \circ b_r \circ c_r \tag{2}$$

where, *R* is a positive integer; $a_r \in \mathbb{R}^I$, $b_r \in \mathbb{R}^J$ and $c_r \in \mathbb{R}^K$ for r=1, ..., R. Elementwise, formula (2) can be written as

$$x_{ijk} \approx \sum_{r=1}^{R} a_{ir} b_{jr} c_{kr}$$

for *i*=1, ..., *I*, *j*=1, ..., *J*, *k*=1, ..., *K* (3)

The Tucker decomposition, a higher-order principal component analysis (PCA), decomposes the target tensor into a core tensor multiplied (or transformed) by a matrix along each mode. In the three-way case where $\mathcal{X} \in \mathbb{R}^{I \times J \times K}$, the Tucker decomposition can be expressed as:

$$\begin{aligned} \mathcal{X} &\approx \mathcal{G} \times_1 A \times_2 B \times_3 C \\ &= \sum_{p=1}^{P} \sum_{q=1}^{Q} \sum_{r=1}^{R} g_{pqr} a_r \circ b_r \circ c_r \end{aligned} \tag{4}$$

where, $A \in \mathbb{R}^{I \times P}$, $B \in \mathbb{R}^{J \times Q}$, and $C \in \mathbb{R}^{K \times R}$ are the factor matrices (usually orthogonal), i.e. the principal components in each mode. The tensor $G \in \mathbb{R}^{P \times Q \times R}$ is called the *core tensor*, whose entries reflect the level of interaction between different components. Elementwise, formula (4) can be written as:

$$x_{ijk} \approx \sum_{r=1}^{R} \sum_{q=1}^{Q} \sum_{r=1}^{R} g_{pqr} a_{ip} b_{jq} c_{kr}$$

for *i*=1, ..., *I*, *j*=1, ..., *J*, *k*=1, ..., *K* (5)

where, P, Q, and R are the number of components (columns) in factor matrices A, B and C, respectively.

The Tensor-Train decomposition represents a *d*-way tensor \mathcal{X} as *d*3-way tensors $\mathcal{X}^{(1)}, \dots, \mathcal{X}^{(d)}$, which are called the TT-cores. The *k*-th TT-core has dimensions r_{k-1} , n_k , r_k where r_{k-1} , r_k are called the TT-ranks. Specifically, each entry of $\mathcal{X} \in \mathbb{R}^{l_1 \times \dots \times l_d}$ can be determined by:

$$a_{i_1\cdots i_d} = \mathcal{X}_{i_1}^{(1)}\cdots \mathcal{X}_{i_d}^{(d)} \tag{6}$$

Next, several important concepts were presented with the following definitions:

Definition 2.1 ([32]) Let $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$ be an *N*-order tensor. The *k*-th mode matrix unfolding is defined as the matrix $X_{(k)} \in \mathbb{R}^{I_k \times J}$, $J = \prod_{i=1, i \neq k}^N I_i$. The tensor element $(i_1, ..., i_N)$ is mapped to the matrix element (i_k, j) , where

$$j = 1 + \sum_{m=1, m \neq k}^{N} (i_m - 1) J_i,$$

with $J_i = \prod_{m=1, m \neq i}^{i-1} I_m.$ (7)

Definition 2.2 ([32]) Let $X \in \mathbb{R}^{I_k \times J}$ be a matrix with $J = \prod_{i=1, i \neq k}^{N} I_i$. The *k*-mode $(I_1, ..., I_N)$ tensor folding \mathcal{X} is defined as the tensor $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$. The matrix element (i_k, j) is mapped to the tensor element $(i_1, ..., i_N)$, where *j* is defined as in formula (7).

Definition 2.3 ([33]) The higher-order singular value decomposition (SVD) of a tensor $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$ can be defined as

$$\mathcal{X} = \mathcal{S} \times_1 U_1 \times_2 U_2 \times_3 \cdots \times_N U_N \tag{8}$$

or elementwise as

$$\begin{aligned} \mathcal{X}(i_1,\cdots,i_N) &= \sum_{j_1=1}^{I_1} \cdots \sum_{j_N=1}^{I_N} \mathcal{S}\left(j_1,\cdots,j_N\right) \cdot \\ & U_1(i_1,j_1) \cdots U_N(i_N,j_N) \end{aligned} \tag{9}$$

where, for $1 \le k \le n$, the U_k are unitary $n_k \times n_k$ matrices, the core tensor $S \in \mathbb{R}^{I_1 \times \cdots \times I_N}$ is such that the (N-1)-order subtensor $S^{i_k = p}$ defined by

$$\mathcal{S}^{l_k=p}_{i_1\cdots i_{k-1}i_{k+1}\cdots i_N} \coloneqq S_{i_1\cdots i_{k-1}pi_{k+1}\cdots i_N} \tag{10}$$

satisfies:

(1) all-orthogonality:

$$\langle \mathcal{S}^{i_k=p}, \mathcal{S}^{i_k=q} \rangle = 0, \forall p \neq q, 1 \le k \le N$$
(11)

(2) the subtensors of the core tensor S are ordered by their Frobenius norm, i.e., for any $1 \le k \le n$,

$$\left\|S^{i_k=1}\right\|_F \ge \dots \ge \left\|S^{i_k=I_N}\right\|_F \ge 0 \tag{12}$$

N-rank of a tensor is the straightforward generalization of the column (row) rank for matrices. Let r_k denote the *k*-th rank of an N-order tensor $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$. Then, it is the rank of the mode-k unfolding matrix $X_{(k)}$:

$$r_k = rank_k(\mathcal{X}) = rank(X_{(k)}) \tag{13}$$

A tensor of which the n-rank are equal to r_n is called a rank-(r_l , ..., r_N) tensor. For any tensor $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$, $vec(\mathcal{X})$ denotes the vectorization of \mathcal{X} .

Next is to introduce the SVD of a matrix. Let there be an $n_1 \times n_2$ matrix of rank r, and $X = U(\text{Diag}\sigma(X))V^T$ be the SVD of X, where U and V are respectively $n_1 \times n$ and $n_2 \times n$ matrices with orthonormal columns, and the function $\sigma: C^{n_1 \times n_2} \rightarrow R^n$ (n = min $\{n_1, n_2\}$) has components $\sigma_1(X) \ge \sigma_2(X) \ge \cdots \ge \sigma_n(X) \ge 0$, i.e. the singular values of the matrix X.

The necessary results in Lewis's research [34] will be briefly described for subsequent formula derivation.

Definition 2.4 ([34]) For any vector γ in \mathbb{R}^n , vector $\tilde{\gamma}$ can be written with components $|\gamma_i|$ arranged in nonincreasing order.

A function *f*: $\mathbb{R}^n \to [-\infty, +\infty]$ is absolutely symmetric, if $f(\gamma) = f(\tilde{\gamma})$ for any vector γ in \mathbb{R}^n .

Theorem 2.1 ([34, Corollary 2.5]) Suppose function f: $R^n \rightarrow (-\infty, +\infty]$ is absolutely symmetric, and $n_1 \times n_2$ matrix X has $\sigma(X)$ in *domf*. Then, the $n_1 \times n_2$ matrix Y lies in $\partial(f \circ \sigma)(X)$, if and only if $\sigma(Y)$ falls in $\partial f(\sigma(X))$ and there exists a simultaneous SVD of the form $U(\text{Diag}\sigma(X))V^T$ and $Y=U(\text{Diag}\sigma(Y))V^T$, where U and V are $n_1 \times n$ and $n_2 \times n$ matrices with orthonormal columns, respectively. In fact,

$$\frac{\partial (f \circ \sigma)(X)}{\partial (f \circ \sigma)(X)} = \{U(\text{Diag}\mu)V^T | \mu \in \partial f(\sigma(x)), X = U(\text{Diag}\sigma(X))V^T\}$$
(14)

2.2 Low n-rank tensor completion problem

This subsection briefly illustrates the low n-rank tensor completion problem. Suppose there exists a partially observed tensor \mathcal{T} under the low n-rank assumption. Then, the Low n-rank tensor completion problem can be expressed as:

$$\min_{\mathcal{X}} \sum_{i=1}^{N} rank(X_{(i)}) \text{ s.t. } \mathcal{X}_{\Omega} = \mathcal{T}_{\Omega}$$
(15)

where, $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$ is the decision variable whose size is identical with \mathcal{T} ; Ω is a subset of index, indicating that the entries of \mathcal{T} in the set Ω are given while the remaining entries are missing. This problem is the special case of the following tensor n-rank minimization problem:

$$\min_{\boldsymbol{\gamma}} \sum_{i=1}^{N} \operatorname{rank}(X_{(i)}) \text{ s.t. } \mathbb{A}(\boldsymbol{\mathcal{X}}) = \boldsymbol{b}$$
(16)

where, the linear map $\mathbb{A}: \mathbb{R}^{I_1 \times \cdots \times I_N} \longrightarrow \mathbb{R}^p$ with $p \leq \prod_{i=1}^N n_i$, and vector $\mathbf{b} \in \mathbb{R}^p$ is given. The rank minimization problem is NP-hard and computationally intractable ([10]). For low-rank matrix minimization problem, the rank function is often replaced with the nuclear norm of a matrix. As an elegant extension of the matrix case, many scholars have attempted to identify a possible convex relaxation as an alternative to the rank minimization problem. For example, Liu et al. proposed the sum of metricized nuclear norm of a tensor:

$$\min_{\mathcal{X}} \sum_{i=1}^{N} \alpha_i \left\| X_{(i)} \right\|_* \text{ s.t. } \mathbb{A}(\mathcal{X}) = \boldsymbol{b}$$
(17)

where, $\alpha_i \ge 0$ (*i*=1, ..., *N*) are constants satisfying $\sum_{i=1}^{N} \alpha_i = 1$; $||Y||_* = \sum_{i=1}^{n} \sigma_i(Y)$ is the nuclear norm of matrix *Y*; $\sigma_i(Y)$ is the i-th largest singular value of matrix *Y*.

In previous work, the model (17) has successfully imputed missing tensor data. Recent studies suggest that the results can be significantly improved, using certain nonconvex functions [35-37].

2.3 DC programming and DC algorithm (DCA)

This subsection briefly outlies DC programming and DCA, facilitating subsequent model solving by DC programming.

Definition 2.4 [38] Let *C* be a convex subset of \mathbb{R}^n . A real-valued function *f*: $C \rightarrow \mathbb{R}$ is called DC on *C*, if there exist two convex functions *g*,*h*: $C \rightarrow \mathbb{R}$ such that *f* can be expressed as:

$$f(x) = g(x) - h(x)$$
 (18)

Definition 2.5 [38] For an arbitrary function $g: R^n \rightarrow R \cup \{+\infty\}$, the function $g^*: R^n \rightarrow R \cup \{+\infty\}$, defined by

$$g^*(y) \coloneqq \sup\{\langle x, y \rangle - g(x) \colon x \in \mathbb{R}^n\}$$
(19)

is called the conjugate function of g.

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Let domg={ $x \in R^n: g(x) < \infty$ } denote the effective domain of *g*. Suppose $x_0 \in \text{dom}g$, the subdifferential of *g* at x_0 , denoted by $\partial g(x_0)$, can be expressed as: $\partial g(x_0):=\{y \in R^n: g(x) \ge g(x_0) + \langle x - x_0, y \rangle, \forall x \in R^n\}$.

As a closed convex set in \mathbb{R}^n , the subdifferential $\partial g(x_0)$ generalizes the derivative in the sense that g is differentiable at x_0 , if and only if $\partial g(x_0)$ is reduced to a singleton which is exactly $\{g'(x_0)\}$.

DC programming is a programming problem dealing with DC functions. The DCA is a continuous approach based on local optimality and the duality of DC programming. The DCA aims to solve the DC program:

$$\beta_p = \inf\{f(x) = g(x) - h(x) : x \in \mathbb{R}^n\}$$
(20)

where, g and h are lower semi-continuous proper convex functions on \mathbb{R}^n . The dual program of formula (20) can be described as:

$$\beta_d = \inf\{h^*(y) - g^*(y) : y \in \mathbb{R}^n\}$$
(21)

Two sequences $\{x_k\}$ and $\{y_k\}$ are optimal solutions of primal and dual programs, respectively. The DCA consists of the two sequences such that the sequences $\{g(x^k)-h(x^k)\}$ and $\{h^*(y^k)-g^*(y^k)\}$ are decreasing. The two sequences are generated as follows: x^{k+1} (resp. y^k) is a solution to the convex program (22) (resp. (23)) defined by

$$\inf\{g(x) - h(x^k) - \langle x - x^k, y^k \rangle \colon x \in \mathbb{R}^n\}$$
(22)

$$\inf\{h^*(y) - g^*(y^{k-1}) - \langle y - y^{k-1}, x^k \rangle: y \in \mathbb{R}^n\}$$
(23)

Problem (22) is obtained from primal DC program (20) by replacing *h* with its affine minimization defined by $h_k(x):=h(x^k)+\langle x-x^k, y^k \rangle$ at a neighborhood of x^k . The solution set of problem (22) is $\partial g^*(y^k)$. Similarly, problem (23) is obtained from the dual DC program (21) through the affine minimization of g^* defined by $(g^*)_k(y):=g^*(y^k)+\langle y-y^k,x^{k+1}\rangle$ at a neighborhood of y^k , and the solution set of problem (23) is $\partial h(x^{k+1})$. Then, the DCA yields the next scheme:

$$y^k \in \partial h(x^k); \ x^{k+1} \in \partial g^*(y^k).$$
(24)

The complete introduction of DC programming and DCA is provided in [38-41].

3. OUR TENSOR COMPLETION ALGORITHM

This section introduces a new model based on the log function, and details our Log-TC algorithm. Our main problem is the following minimization problem:

min
$$\Phi(\mathcal{X})$$
 s.t. $\mathcal{A}(\mathcal{X}) = \mathbf{b}$ (25)

where, $\Phi(\mathcal{X}) = \sum_{i=1}^{N} \varphi(X_{(i)})$, $\varphi(X) = \sum_{j=1}^{n} \log (\sigma_j(X_{(i)}) + \varepsilon)$. In essence, the log function acting on a tensor *X* is a convex combination of the log functions acting on all matrices $X_{(i)}$ unfolded along each mode. Specifically, the log function acting on the matrix $X_{(i)}$ is actually a convex combination of the log function of the log function acting on the singular value of this matrix, i.e.

 $\sum_{j=1}^{n} \log (\sigma_j(X_{(i)}) + \varepsilon)$. Next, an unconstrained problem was considered:

$$\min \Phi(\mathcal{X}) + \frac{1}{2\mu} \|\mathcal{A}(\mathcal{X}) - \mathbf{b}\|_2^2$$
(26)

where, $\mu > 0$ is a penalty parameter.

The distributions of rank function, l_1 function, and log function were compared to further explain the different approximation effects.



Figure 1. Comparison of rank function, *l*₁ function, and log function

As shown in Figure 1, log function fell between rank function and l_1 norm, which can simultaneously increase the penalty on small values and decrease the penalty on large values.

Next, the DCA mentioned in Subsection 2.3 was employed to solve model (26).

Let $f(\mathcal{X}) = \Phi(\mathcal{X}) + \frac{1}{2\mu} \|\mathcal{A}(\mathcal{X}) - \mathbf{b}\|_2^2$. Then model (26) can be expressed as:

$$\min\{f(\mathcal{X}) := g(\mathcal{X}) - h(\mathcal{X})\}$$
(27)

where, $\mathcal{X} \in \mathbb{R}^{n_1 \times n_2 \times \cdots n_N}$. The DC components g(X) and h(X) can be respectively given by:

$$g(\mathcal{X}) = \frac{1}{2\mu} \|\mathcal{A}(\mathcal{X}) - \mathbf{b}\|_2^2 + \|\mathcal{X}\|_*$$
(28)

where, $\lambda \ge 0$ is a parameter; and

$$h(\mathcal{X}) = \|\mathcal{X}\|_* - \Phi(\mathcal{X}) \tag{29}$$

According to Subsection 2.3, the DCA scheme for model (27) was determined by computing the two sequences $\{X^k\}$ and $\{Y^k\}$ which satisfy the following conditions:

$$\mathcal{Y}^{k} \in \partial h(\mathcal{X}^{k}), \, \mathcal{X}^{k+1} \in \partial g^{*}(\mathcal{Y}^{k})$$
(30)

The calculation process of $\partial h(X^k)$ and $\partial g^*(Y^k)$ will be repeated in the following subsections.

3.1 Calculation of $\partial h(X^k)$

From the previous discussion, we have:

$$h(\mathcal{X}^{k}) = \|\mathcal{X}^{k}\|_{*} - \Phi(\mathcal{X}^{k}) = \sum_{i=1}^{N} (\alpha_{i} \|X_{(i)}^{k}\|_{*} - \varphi(X_{(i)}^{k}))$$
(31)

where, $\Phi(\mathcal{X}) = \sum_{i=1}^{N} \varphi(X_{(i)})$; $\varphi(X) = \sum_{j=1}^{n} \log (\sigma_j(X_{(i)}) + \varepsilon)$. Then, the subdifferential of $h(X^k)$ is:

$$\frac{\partial h(\mathcal{X}^k) = \sum_{i=1}^N \partial(\alpha_i \|X_{(i)}^k\|_* - \varphi(X_{(i)}^k)) =}{\sum_{i=1}^N (\alpha_i \partial f_1 \circ \sigma(X_{(i)}^k) - \partial f_2 \circ \sigma(X_{(i)}^k))}$$
(32)

where, $\sigma(X)$ denotes is the vector consisting of the singular values of matrix X, and $\forall x \in \mathbb{R}^n$: $f_1(x) = \sum_{i=1}^n x_i$ and $f_2(x) = \sum_{i=1}^n \log(x_i + \epsilon)$.

By Theorem 2.1, we have:

$$Y \in \partial(f \circ \sigma)(X) = \{U(\text{Diag}\omega)V^T | \omega \in \partial f(\sigma(X)), X = U(\text{Diag}\sigma(X))V^T\}$$
(33)

Thereby, following conclusion can be drawn:

$$U\begin{pmatrix} \alpha_{i} & 0 & 0 & 0\\ 0 & \alpha_{i} & 0 & 0\\ 0 & 0 & \ddots & 0\\ 0 & 0 & 0 & \alpha_{i} \end{pmatrix}_{r_{i} \times r_{i}} V^{T} \in \alpha_{i} \partial f_{1} \circ \sigma(X_{(i)}^{k})$$
(34)

and

$$U\begin{pmatrix} \frac{1}{\sigma_{1}(X_{(i)}^{k})+\varepsilon} & 0 & 0 & 0\\ 0 & \frac{1}{\sigma_{2}(X_{(i)}^{k})+\varepsilon} & 0 & 0\\ 0 & 0 & \ddots & 0\\ 0 & 0 & 0 & \frac{1}{\sigma_{r_{i}}(X_{(i)}^{k})+\varepsilon} \end{pmatrix}_{r_{i}\times r_{i}}$$
(35)
$$V^{T} \in \partial f_{2} \circ \sigma(X_{(i)}^{k})$$

where, r_i is the rank of matrix $X_{(i)}^k$. Therefore, an element Y^k could be found in $\partial h(X^k)$ which has the following form:

$$\mathcal{Y}^k = \frac{1}{N} \sum_{i=1}^N foldM_i \tag{36}$$

where,

$$M_{i} = U \begin{pmatrix} \alpha_{i} - \frac{1}{\sigma_{i}(x_{(i)}^{k}) + \varepsilon} & 0 & 0 & 0 \\ 0 & \alpha_{i} - \frac{1}{\sigma_{z}(x_{(i)}^{k}) + \varepsilon} & 0 & 0 \\ 0 & 0 & \ddots & 0 \\ 0 & 0 & 0 & \alpha_{i} - \frac{1}{\sigma_{r_{i}}(x_{(i)}^{k}) + \varepsilon} \end{pmatrix} V^{T} \quad (37)$$

3.2 Calculation of $\partial g^*(Y^k)$

According to the preliminary knowledge in Subsection 2.3, $\partial g^*(Y^k)$ is a solution to the convex program (22). $\forall \mathcal{Y}^k \in \mathbb{R}^{n_1 \times n_2 \times \cdots \cdot n_N}$,

$$\partial g^*(\mathcal{Y}^k) = \arg \min\{g(\mathcal{X}) - h(\mathcal{X}^k) - \langle \mathcal{X} - \mathcal{X}^k, \mathcal{Y}^k \rangle \\ : \mathcal{X} \in \mathbb{R}^{n_1 \times n_2 \times \cdots n_N}\}$$

$$= \arg \min\{g(\mathcal{X}) - \langle \mathcal{X} - \mathcal{X}^k, \mathcal{Y}^k \rangle : \mathcal{X} \in \mathbb{R}^{n_1 \times n_2 \times \cdots n_N}\}$$

$$(38)$$

Combining (28) and (38), the following optimization can be obtained:

$$\min \frac{1}{2\mu} \|\mathcal{A}(\mathcal{X}) - \mathbf{b}\|_2^2 + \|\mathcal{X}\|_* - \langle \mathcal{X} - \mathcal{X}^k, \mathcal{Y}^k \rangle$$

$$, \mathcal{Y}^k \in \partial h(\mathcal{X}^k)$$
(39)

Since the optimal condition for minimization of a convex function is $0 \in \partial \varphi(x)$, \mathcal{X}^* is optimal of formula (39) if and only if

$$0 \in \frac{1}{\mu} \mathbb{P}(\mathcal{X}^*) + \partial \|\mathcal{X}\|_* - \mathcal{Y}^k \tag{40}$$

where, $P(\mathcal{X}^*) = \mathcal{A}^* (\mathcal{A}(\mathcal{X}^*) - b)$ and \mathcal{A}^* represents the adjoint of \mathcal{A} .

For $\forall \tau > 0$, formula (40) is equivalent to:

$$0 \in \frac{1}{\mu}\tau P(\mathcal{X}^*) + \tau \,\partial \|\mathcal{X}\|_* - \tau \mathcal{Y}^k + \frac{1}{\mu}\mathcal{X}^* - \frac{1}{\mu}\mathcal{X}^* \tag{41}$$

Then, it holds that

$$0 \in \tau \,\partial \|\mathcal{X}\|_* + \frac{1}{\mu}\mathcal{X}^* - \frac{1}{\mu}\big(\mathcal{X}^* - \tau \mathsf{P}(\mathcal{X}^*)\big) - \tau \mathcal{Y}^k \tag{42}$$

Let $\mathcal{Z}^* = \mathcal{X}^* - \tau P(\mathcal{X}^*)$. Formula (42) can be reduced to:

$$0 \in \tau \,\partial \|\mathcal{X}\|_* + \frac{1}{\mu}\mathcal{X}^* - \frac{1}{\mu}\mathcal{Z}^* - \tau \mathcal{Y}^k \tag{43}$$

After combining, we have:

$$0 \in \tau \,\partial \|\mathcal{X}\|_* + \frac{1}{\mu}(\mathcal{X}^* - (\mathcal{Z}^* + \mu\tau\mathcal{Y}^k)) \tag{44}$$

i.e. \mathcal{X}^* is the optimal solution to

$$\min \tau \|\mathcal{X}\|_* + \frac{1}{2\mu} \|\mathcal{X} - (\mathcal{Z}^* + \mu \tau \mathcal{Y}^k)\|_F^2$$
(45)

Under the definition of mode-n unfolding, the problem of minimizing (45) can be written as:

$$\min \tau \sum_{i=1}^{N} \alpha_i \|X_{(i)}\|_* + \frac{1}{2N\mu} \sum_{i=1}^{N} \|X_{(i)} - (Z_{(i)}^* + \mu\tau Y_{(i)}^k)\|_F^2$$
(46)

Problem (46) is difficult to solve due to the interdependent nuclear norms. Therefore, a series of new matrix variables M_i (*i*=1, 2, ..., *N*) were introduced such that they equal to $X_{(1)}$, $X_{(2)}, ..., X_{(N)}$, respectively, which represent the different moden unfoldings of the tensor *X*. With these new variables, formula (46) can be rewritten as:

$$\min \tau \sum_{i=1}^{N} \alpha_i \|M_i\|_* + \frac{1}{2N\mu} \sum_{i=1}^{N} \|M_i - (Z_{(i)}^* + \mu\tau Y_{(i)}^k)\|_F^2 \text{ s.t. } M_i = X_{(i)}, \forall i \in \{1, 2, \cdots, N\}$$
(47)

Relaxing the constraint $M_i=X_{(i)}$, the following problem can be obtained:

$$\min \tau \sum_{i=1}^{N} \alpha_{i} \|M_{i}\|_{*} + \frac{1}{2N\mu} \sum_{i=1}^{N} \|M_{i} - (Z_{(i)}^{*} + \mu\tau Y_{(i)}^{k})\|_{F}^{2} + \frac{1}{2\beta} \sum_{i=1}^{N} \|M_{i} - X_{(i)}\|_{F}^{2}$$
(48)

where, β >0 is a penalty parameter. An optimal solution of problem (48) approaches an optimal solution of formula (47) as β >0 [42]. For convenience, suppose $\beta = \mu$ in problem (48) and consider the following minimization:

$$\min_{\mathcal{X},M_{i}} \tau \sum_{i=1}^{N} \alpha_{i} \|M_{i}\|_{*} + \frac{1}{2N\mu} \sum_{i=1}^{N} \|M_{i} - (\mathbf{Z}_{(i)}^{*} + \mu\tau \mathbf{Y}_{(i)}^{k})\|_{F}^{2} + \frac{1}{2\mu} \sum_{i=1}^{N} \|M_{i} - X_{(i)}\|_{F}^{2}$$
(49)

Computing M_i Fixing all variables except M_i (*i*=1, 2, ..., *N*), problem (49) can be transformed into the following matrix problem:

$$\min_{M_{i}} \tau \alpha_{i} \|M_{i}\|_{*} + \frac{1}{2N\mu} \|M_{i} - (Z_{(i)}^{*} + \mu \tau Y_{(i)}^{k})\|_{F}^{2} + \frac{1}{2\mu} \|M_{i} - X_{(i)}\|_{F}^{2}$$
(50)

The optimal solution of problem (50) will be searched for by the following theorem.

Theorem 3.1 Suppose $\mu > 0$ and $\tau > 0$. For any given $i \in \{1, 2, ..., N\}$, M_i^* is an optimal solution to problem (50) if and only if

$$M_i^* = D_{\frac{\tau \alpha_i \mu N}{1+N}} \left(\frac{Z_{(i)}^* + \mu \tau Y_{(i)}^k + N X_{(i)}}{1+N} \right)$$
(51)

Proof: M_i^* is an optimal solution to problem (50) if and only if

$$0 \in \partial \tau \alpha_{i} \|M_{i}^{*}\|_{*} + \frac{1}{N\mu} (M_{i}^{*} - (Z_{(i)}^{*} + \mu \tau Y_{(i)}^{k})) + \frac{1}{\mu} (M_{i}^{*} - X_{(i)})$$
(52)

where, $\partial \|M_i^*\|_*$ is the subgradients of $\|\cdot\|_*$ at M_i^* . Through proper combining and deforming, we have:

$$0 \in \partial \frac{\tau \alpha_{i}}{\frac{1+N}{\mu N}} \|M_{i}^{*}\|_{*} + M_{i}^{*} - \frac{\frac{1}{\mu N}}{\frac{1+N}{\mu N}} (Z_{(i)}^{*} + \mu \tau Y_{(i)}^{k}) - \frac{\frac{1}{\frac{\mu}{1+N}} X_{(i)}}{\frac{1}{\frac{1+N}{\mu N}} X_{(i)}}$$
(53)

which is equivalent to

$$0 \in \partial \frac{\tau \alpha_i \mu N}{1+N} \| M_i^* \|_* + M_i^* - \frac{1}{1+N} (Z_{(i)}^* + \mu \tau Y_{(i)}^k + N X_{(i)})$$
(54)

i.e. M_i^* is an optimal solution to the following minimization problem

$$\min \frac{\tau \alpha_{i\mu N}}{1+N} \|M_{i}\|_{*} + \frac{1}{2} \left\| M_{i} - \frac{1}{1+N} \left(Z_{(i)}^{*} + \mu \tau Y_{(i)}^{k} + N X_{(i)} \right) \right\|_{F}^{2}$$
(55)

According to the Theorem 3 in Ma et al. [43], an optimal solution to problem (55) must be the matrix shrinkage operator applied to $\frac{1}{1+N} (Z_{(i)}^* + \mu \tau Y_{(i)}^k + NX_{(i)})$. Hence, the optimal solution of problem (55) is

$$M_{i}^{*} = D_{\frac{\tau \alpha_{i} \mu N}{1+N}} \left(\frac{Z_{(i)}^{*} + \mu \tau Y_{(i)}^{k} + N X_{(i)}}{1+N} \right)$$
(56)

For $\forall v \in \mathbb{R}^+$, the matrix shrinkage operator $D_v(X)$ can be defined as:

$$D_{\nu}(X) = UDiag(\bar{\sigma}(X))V^{T}$$
(57)

Let $X = UDiag(\sigma(X))V^T$ be the SVD for matrix X,

$$\bar{\sigma}(X) = \begin{cases} \sigma_i(X) - \nu, & \text{if } \sigma_i(X) - \nu > 0\\ 0, & o.w. \end{cases}$$
(58)

Then, solution (56) is also the optimal solution to problem (50).

Q.E.D.

Computing *X*: Fixing all other variables except $X_{(i)}$, *i*=1, 2, ..., *N*, the following problem was considered to get the exact solution:

$$\min \frac{1}{2\mu} \|M_i - X_{(i)}\|_F^2 \text{ s.t. } X_{(i)}^* = M_i$$
(59)

From the previous discussion, M_i can be regarded as known in model (59). Hence, the solution to model (59) can be obtained easily as:

$$\mathcal{X}^* = \frac{1}{N} \sum_{i=1}^{N} fold(M_i)$$
(60)

From formulas (56) and (60), the following iterative formula can be derived:

$$\begin{pmatrix} \alpha_{i} - \frac{1}{\sigma_{1}(x_{(i)}^{k}) + \varepsilon} & 0 & 0 & 0 \\ 0 & \alpha_{i} - \frac{1}{\sigma_{2}(x_{(i)}^{k}) + \varepsilon} & 0 & 0 \\ 0 & 0 & \ddots & 0 \\ 0 & 0 & 0 & \alpha_{i} - \frac{1}{\sigma_{r_{i}}(x_{(i)}^{k}) + \varepsilon} \end{pmatrix} V^{T} \\ Z_{(i)}^{k} = \mathcal{X}^{k} - \tau P(\mathcal{X}^{k}) \\ \mathcal{X}^{k+1} = \frac{1}{N} \sum_{i=1}^{N} fold \ D_{\frac{\tau \alpha_{i} \mu N}{1+N}} \left(\frac{Z_{(i)}^{k} + \mu \tau Y_{(i)}^{k} + N X_{(i)}}{1+N} \right)$$

$$(61)$$

Through the above discussion, the Log-TC algorithm can be established to solve tensor completion problem (Algorithm 1).

Algorithm 1: Log-TC algorithm
Input: \mathcal{A} , b, μ , and τ
Initialization: Set k:=0, and $\mathcal{X}^k \coloneqq 0$
for $\mu \rightarrow 0$, do
while not converged, do
$P(\mathcal{X}^k) = \mathcal{A}^*(\mathcal{A}(\mathcal{X}^*) - b)$
$\mathcal{Z}^k = \mathcal{X}^k - \tau P(\mathcal{X}^k)$
for $i = 1:N$
$M_{i}^{k} = D_{\frac{\tau \alpha_{i} \mu N}{1+N}} \left(\frac{Z_{(i)}^{k} + \mu \tau Y_{(i)}^{k} + N X_{(i)}}{1+N} \right)$
end
$\mathcal{X}^{k+1} = \frac{1}{N} \sum_{i=1}^{N} fold(M_i^k)$
end while
end for
Output: X_{ult}

4. NUMERICAL EXPERIMENTS

To verify its empirical performance, the proposed Log-TC algorithm was compared with fixed point iterative method for low-rank tensor completion (FP-LRTC) [44] and fast low rank tensor completion algorithm (FaLRTC) [10] through numerical experiments. The FP-LRTC, which combines a fixed-point iterative method for solving the low n-rank tensor pursuit problem, and a continuation technique, has been applied to solve the relaxation model of the low-rank tensor completion problem. The FaLRTC utilizes a smoothing scheme to transform the original nonsmooth problem into a smooth problem, and was found more efficient than simple low rank tensor completion (SiLRTC), and high accuracy low rank tensor completion (HaLRTC).

This section is separated into two subsections. In Subsection 4.1, our algorithm was tested on the randomly generated low n-rank tensors. In Subsection 4.2, our algorithm was tested on

color image recovery problem. All experiments were carried out in MATLAB R2016a on a computer running on Windows 10, with a 3.20GHz CPU and 4GB of memory. For fair comparison, each number is the average over 10 separate runs.

4.1 Synthetic data

First, our Log-TC algorithm was tested on several synthetic datasets for the tensor completion tasks. Following the method proposed by Rauhut et al. [45], a n-order tensor $\mathcal{T} \in \mathbb{R}^{n_1 \times n_2 \times \cdots \times n_N}$ was created randomly with the *n*-rank (r_1, r_2, \ldots, r_N) of the form $\mathcal{T} \coloneqq \mathcal{M} \times_1 U_1 \times_2 \cdots \times_N U_N$, where $\mathcal{M} \in \mathbb{R}^{r_1 \times r_2 \times \cdots \times r_N}$ is the core tensor and $U_i \in \mathbb{R}^{n_i \times r_i} (i = 1, 2, \cdots, N)$. Each entry of the core tensor \mathcal{M} was sampled independently from a standard Gaussian distribution $\mathcal{N}(0, 1)$. Then, m entries were sampled randomly from tensor \mathcal{T} in a uniform manner. The sample ratio, denoted by *SR*, was defined as $SR = m/I_1 \cdots I_N$. The relative error (RelErr) of the recovered tensor defined as:

$$\operatorname{RelErr} = \frac{\|x_{ult} - \mathcal{T}\|_F}{\|\mathcal{T}\|_F}$$
(62)

is used to estimate the closeness of \mathcal{X}_{ult} to \mathcal{T} , where \mathcal{X}_{ult} is the optimal solution to model (26) produced by the algorithms.

The ratio "fr" [44] between the degree of freedom (DOF) in a rank r matrix to the number of samples in the matrix completion was extended to tensor completion problem and denoted by FR:

$$FR = \frac{1}{N} \sum_{i=1}^{N} fr_i = \frac{1}{N} \sum_{i=1}^{N} \frac{r_i(I_i + \Pi_i - r_i)}{I_1 \cdots I_N \times SR}$$
(63)

where, $r_i = \operatorname{rank}(X_{(i)})$; $\Pi_i = \prod_{k=1, k \neq i}^N I_k$. In the following results, "FR" is used to represent the hardness of the tensor completion. If FR is greater than 0.6, then the problems are "hard problems"; otherwise, the problems are "easy problems". The parameters wee configured as: $\mu = 1$, $\theta_{\mu} = 1$ -sr, $\bar{\mu} = 1 \times 10^{-8}$, $\tau = 10$, and $\lambda = 1$.

The simulated performance of FP-LRTC, FaLRTC and

Log-TC for random tensor completion problem, as measured by relative error and running time, are reported in Figure 2 and Table 1.

As shown in Figure 2, the relative error for tensor of size $60 \times 60 \times 60$ with the n-rank was fixed at (6,6,6), and the sample rate changed from 10% to 80%. As it is known, when the n-rank is fixed, the complexity of the problem increases with the decrease of the SR. The plot illustrates that Log-TC is clearly superior to FP-LRTC and FaLRTC. The superiority is particularly evidence in the situation of SR=10%: Log-TC recovered the tensor with the relative error about 10⁻⁴, while FaLRTC and FP-LRTC with 10⁰.



Figure 2. The relative error on tensors of size $60 \times 60 \times 60$ with sample rate (SR) between 10% to 80% and fixed n-rank (6,6,6)

Table 1. Comparison of running time on tensors of size $60 \times 60 \times 60$ with sample rate (SR) 70% and -rank (6,6,6)

Relative	Log-TC	FP-LRTC	FaLRTC
Error	Time (s)	Time (s)	Time (s)
10-9	67.23	79.04	132.41
10^{-8}	26.62	32.75	40.42
10^{-6}	11.07	16.36	36.21
10^{-4}	2.82	4.51	41.17

Table 1 shows the performance of each algorithm with SR 70% in terms of the running time. Log-TC obviously outperformed the contrastive algorithms, under the same relative error. Even when the precision requirement is ultrahigh (10⁻⁹), Log-TC can complete the trask in one minute.

Figure 3 shows the relative error for tensor of size $60 \times 60 \times 60$, when the sample rate was fixed at SR=40% and n-

rank r changed from 2 to 40. Like the previous situation, when sample rate is fixed, the complexity of the problem increases with the increase of the rank. As shown in Figure 3, Log-TC outperformed FP-LRTC and FaLRTC in all cases. Log-TC performed especially well at r=20: Log-TC recovered the tensors with the relative error about 10^{-8} , while FP-LRTC and FaLRTC with only about 10^{-1} . It can also be seen that, when r=24, our algorithm solved the problem with the relative error 10^{-3} , but the other two methods failed to effectively recover the tensors.



Figure 3. The relative error on tensors of size $60 \times 60 \times 60$ n-rank (r,r,r) between 2 and 40 and fixed sample rate (SR=40%).

Next, the numerical results of the three algorithms on easy problems and hard problems were reported. Tables 2 and 3 list the numerical results of several cases in which the test tensors have different sizes, n-ranks and sample rates.

According to the results on easy problems (Table 2), the Log-TC outshined the other two algorithms in relative error and running time. FaLRTC is inferior to FP-LRTC and Log-TC in terms of relative error. FP-LRTC and Log-TC performed similarly on relative error, but Log-TC consumed the shorter time. Overall, Log-TC achieved the best performance on easy problems.

According to the results on hard problems (Table 3), Log-TC also realized better performance than the contrastive algorithms. With the increase of tensor order, the robustness of FaLRTC was getting worse. FP-LRTC and Log-TC were always capable of recovering the tensors efficiently, but the running time of FP-LRTC is about 2.5 times more than that of the Log-TC.

Table 2. Comparison results for easy problems

Tensor	Rank	сD	FR	Log-TC		FP-LRTC		FaLRTC	
		SR		RelErr	Time	RelErr	Time	RelErr	Time
20×30×40	(2,2,2)	0.5	0.15	1.72e-08	1.11	3.36e-08	1.82	1.23e-06	3.53
30×30×30	(5,5,5)	0.3	0.57	5.42e-08	3.24	8.27e-07	4.39	4.83e-06	2.72
60×60×60	(5,5,5)	0.7	0.12	4.57e-09	2.83	6.71e-09	4.65	1.11e-05	7.88
100×100×100	(10,10,10)	0.4	0.25	3.95e-09	50.40	5.90e-09	87.28	1.21e-05	36.63
20×20×30×30	(5,5,5,5)	0.6	0.35	4.40e-09	13.57	5.76e-09	22.87	5.94e-07	90.76
50×50×50×50	(6,6,6,6)	0.7	0.17	1.23e-09	90.77	1.64e-09	139.19	2.36e-08	189.6
30×30×30×30×30	(2,2,2,2,2)	0.5	0.13	1.83e-09	635.81	2.44e-09	1573.9	9.83e-08	2896
30×30×30×30×30	(6,6,6,6,6)	0.6	0.33	1.43e-09	905.60	1.47e-09	1649.1	8.63e-08	2765

Table 3. Comparison results for hard problems

Tensor	Rank	SR	FR	Log-TC		FP-LRTC		FaLRTC	
	Kalik			RelErr	Time	RelErr	Time	RelErr	Time
20×30×40	(2,2,2)	0.1	0.75	3.90e-04	50.82	5.72e-01	67.61	2.11e-01	6.73
30×30×30	(5,5,5)	0.1	1.71	8.30e-10	57.83	8.45e-7	62.09	8.74e-01	1.31
70×70×70	(10, 10, 10)	0.2	0.72	2.28e-08	324.83	2.71e-08	422.34	2.35e-04	56.42
100×100×100	(10, 10, 10)	0.2	0.63	1.20e-08	119.29	1.67e-08	187.10	1.09e-04	191.5
20×20×30×30	(5,5,5,5)	0.3	0.7	1.34e-08	241.36	4.76e-07	373.09	3.42e-07	896.3
50×50×50×50	(5,5,5,5)	0.2	0.69	7.18e-09	754.08	5.43e-08	1271.51	2.72e-01	1732
20×20×20×20×20	(5,5,5,5,5)	0.3	0.83	3.04e-08	398.47	6.28e-06	1689.23	3.92e-01	2874
20×20×20×20×20	(8,8,8,8,8)	0.4	1.00	4.48e-09	746.92	8.52e-08	1939.63	4.53e-08	2145

4.2 Image simulation

This subsection compared the three algorithms on the recovery of color images. Color images can be expressed as third order tensors. If the images are of low rank, the image recovery can be solved as a low n-rank tensor completion problem. Assuming that the images are well structured, FP-LRTC, FaLRTC and Log-TC were applied separately to solve the recovery problem.

The recovery effects of the Log-TC are illustrated in Figure 4 and Table 4. In Figure 4, (a1), (b1), (c1), and (d1) are original color images. Then, 80%, 60%, 40% and 20% entries were randomly removed, creating corrupted images (a2), (b2), (c2), and (d2), respectively. The images recovered by the Log-TC are (a3), (b3), (c3), and (d4), respectively. Only the effects of the Log-TC were provided, because the three algorithms achieved similar recovery accuracy. The recovery accuracy and running time of each algorithm are listed in Table 4.



	Log-TC		FP-LR	атс	FaLRTC		
SR	Relative	Time	Relative	Time	Relative	Time	
	Error	(s)	Error	(s)	Error	(s)	
20%	9.14e-02	1300	9.62e-02	1681	8.31e-02	2137	
40%	6.36e-02	598	8.45e-02	740	7.74e-02	1402	
60%	4.89e-02	362	5.71e-02	298	2.57e-02	978	
80%	3.40e-02	209	4.21e-02	302	3.91e-02	721	

As shown in Figure 4, the Log-TC recovered the natural images with missing information. Table 4 shows that our algorithm achieved a recovery accuracy of about 10^{-2} , and a clear edge over the contrastive methods in running time.



Figure 5. Original images, corrupted images, recovered images by the Log-TC in the recovery of images with poor texture



Figure 4. Original images, corrupted images, and recovered image by Log-TC in the recovery of natural images

Table 5. Numerical results on the color image with poor texture

SR	Log-T	С	FP-LRT	C	FaLRTC		
эк	Relative Error Time(Relative Error Time(s)		Relative Error	Time(s)	
20%	1.17e-01	1835	5.72e-01	1920	4.53e-01	2341	
40%	6.71e-02	857	8.45e-02	1103	7.47e-02	1780	
60%	5.02e-02	493	2.71e-02	631	3.87e-02	1285	
80%	3.37e-02	291	4.42e-02	329	2.57e-02	943	

The original images in Figure 4 have a nice mixture of details, shading areas, flat regions, and textures. Next, the three algorithms were separately adopted to recover color images with poor texture. The recovery effect of our algorithm is displayed in Figure 5, and the numerical results of the three methods are compared in Table 5. The subgraphs in the middle column of Figure 5 were prepared in the same method as those in the middle column of Figure 4. It can be seen that, despite the poor texture of images, our algorithm still recovered the details of the images, and outperformed the other algorithms in running time.

5. CONCLUSIONS

This paper firstly defines a log function of tensor, and uses it as an approximation of tensor rank function. Furthermore, a simple yet efficient algorithm named Log-TC algorithm was proposed to recover an underlying low n-rank tensor. The log function strikes a balance between rank minimization and nuclear norm. It can simultaneously increase the penalty on small values and reduce the penalty on large values. After that, our algorithm was compared with FP-LRTC and FaLRTC through numerical experiments on randomly generated tensors and color image recovery. The numerical comparisons demonstrate that the Log-TC outperformed the other two algorithms. The future research will investigate the largescale cases by other techniques, and test the performance of our algorithm for image inpainting problems.

ACKNOWLEDGMENT

The authors would like to thank the anonymous reviewers for the useful comments and suggestions. This work is partially supported by National Natural Science Foundation of China Grant (No. 11626080), Natural Science Foundation of Hebei Province (No. A2017207011), National Statistical Science Research Project of China (No. 2019LY27) and Department of Education of Hebei Province under Grant (No. QN2018018).

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