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#### Research article

# A new algorithm by embedding structured data for low-rank tensor ring completion

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**Abstract:** In this paper, we put up with a new algorithm for tensor completion problems that include missing slices or row/column fibers, where embedding a structured tensor by a multi-way delayembedding transform (MDT) makes the tensor to be completed have a special structure. The main idea is to employ a tensor completion algorithm based on the tensor ring rank, constructing latent tensor ring factors with a structure that approximates the original tensor starting from the tensor structure. It is also proved that the sequence generated by the new algorithm converges to the optimal solution. Finally, the feasibility of the proposed algorithm is verified by experiments. Compared with other completed algorithms based on tensor ring rank, the completed accuracy is improved, up to 30%.

**Keywords:** low-rank tensor completion; tensor ring decomposition; embedded space; structured data **Mathematics Subject Classification:** 15A69, 90C06, 90C47

## 1. Introduction

The problem of completing a low-rank matrix or tensor from small sets of linear measurements occurs in many areas, such as hyperspectral compressive sensing [1], color images [2, 3], data analysis [4], various video processing [5, 6], and machine learning [7]. Mathematically, the optimization model of a tensor completion is described as follows [8]:

$$\min_{\mathcal{X}} \operatorname{rank}(\mathcal{X}),$$
s.t.  $P_{\Omega}(\mathcal{X}) = P_{\Omega}(\mathcal{T}),$  (1.1)

where  $\mathcal{X}, \mathcal{T} \in \mathbb{R}^{I_1 \times I_2 \times ... \times I_N}$  are both Nth-order tensors,  $\mathcal{T}$  is the observed incomplete tensor, and  $\Omega$  denotes the index set of the known samples,  $P_{\Omega}(\cdot) : \mathbb{R}^{I_1 \times I_2 \times ... \times I_N} \to \mathbb{R}^{I_1 \times I_2 \times ... \times I_N}$  is the associated sampling orthogonal projection operator, which acquires only the entries indexed by the set  $\Omega$ .

As we know, the rank of a tensor has many definitions based on different decompositions of a tensor. The most popular ones currently include the CANDECOMP/PARAFAC (CP) rank [9], Tucker rank (or multilinear rank) [10], tensor train (TT) rank [11], and tensor ring (TR) rank [12]. These tensor ranks have found application in various fields. The tensor completion problem is one of the most important applications of tensor decompositions, with the goal to complete an incomplete tensor from the available data. Moreover, the rank-minimization-based tensor completion models can be generated by citing different rank function; it is very challenging to specify the optimal rank beforehand. This is especially the case for Tucker, TT, and TR decompositions, for which the rank is defined as a vector; it is therefore impossible to find the optimal ranks by cross-validation due to the immense possibilities. And many theoretical results and algorithmic methods for completing a low-rank tensor have been obtained. For example, Liu et al. [8] first proposed a definition for the tensor nuclear norm, minimize the weighted sum of the nuclear norm on unfolding matrices to tackle the color image and video completion problems. Then (1.1) can be relaxed into a convex optimization problem as follows:

$$\min \sum_{i=1}^{N} \alpha_i ||X_{(i)}||_*,$$
  
s.t.  $P_{\Omega}(X) = P_{\Omega}(\mathcal{T}),$  (1.2)

where  $\alpha_i \geq 0$  are constants satisfying  $\sum_{i=1}^N \alpha_i = 1$ ,  $\mathcal{X}_{(i)} \in \mathbb{R}^{I_i \times \prod_{j \neq i} I_j}$  is the mode-*i* unfolding of tensor X for  $i \in \{1, ..., N\}$ , and  $||X_{(i)}||_*$  is the nuclear norm of matrix  $X_{(i)}$ . Later,  $X_{(i)}$  and  $X_{(i)}$  are all [13] and Liu et al. [14] estimated the missing data with the regularized CP decomposition. Furthermore, Bengua et al. [15] applied the TT format to recover incomplete tensors. Zhang et al. [16] considered minimizing a tensor nuclear norm (TNN) on t-SVD to find a low tubal-rank tensor. Zhao et al. [17] employed a Bayesian treatment on the CP model to handle the incomplete tensor. Chen et al. [18] further explored a Bayesian Gaussian CP decomposition model on the traffic data imputation problem. Long [19] presented a Bayesian low-rank tensor completion method with the TR format for the image completion problem. The central effort, for which this paper will further study, focuses on how to retain the correlated information of data sets as much as possible under the tensor low-rank framework in a tractable way in order to achieve the high accuracy of approximations. Based on TR decomposition, Yuan et al. [20] introduced the weight optimization algorithm (TR-WOPT) and the singular value thresholding algorithm (TR-SVD); Zhao et al. [21] proposed the least squares method (TR-ALS) to approximate tensors. Recently, some works [22, 23] jointly consider the low-rank and smoothness properties, which use the hierarchical structured information. Besides, the structured low-rank model also proposed in [24]. Numerous detailed derivations on the tensor completion problem can be referred to in the references given therein.

These algorithms mentioned above were usually used to generalize tensor completion models. They directly represent data with the tensors themselves, and some structures of the tensors are assumed, and then they can be considered as one of two classes of methods for modeling tensors. By contrast, the algorithms in the second class are embedding the data into a high-dimensional feature space, and it is assumed that the data can be represented by a low-rank or a smooth manifold in the embedded space [25–28]. Typically, a time series is represented by a Hankel matrix, and its low-rank property

has been employed widely for modeling a linear time-invariant system of singles [27,28]. For example, Li et al. [29] proposed a method for modeling damped sinusoidal signals based on a low-rank Hankel approximation. Ding et al. [25] proposed the use of rank minimization of a Hankel matrix for the video inpainting problem by assuming an autoregressive moving average model.

This study focuses on a case where all of the elements in some continuous slices or row/column fibers are missing in tensor data. The completion of missing slices or row/column fibers is recognized as a difficult problem that common tensor completion algorithms usually fail to solve in this case. The research on using embedded structured data for low rank or smoothness of data has mainly been applied to matrix completion problems; only a few works [30, 31] have studied tensor completion algorithms for embedded structures. Therefore, there is still a lot of research in tensor completion problems for embedding structured data.

To obtain the structured data, the popular method of embedding data is vector delay embedding (VDT) or multi-way delay-embedding transform (MDT) [31], which is a special Hankel structure formed by embedding a set of vectors into some data. It is frequently used in the signal processing of time series [32]. It is noted that [31] have proposed two Tucker decomposition algorithms by using the same embedding method. This paper is to use the MDT technique to embed data from general tensors and expand them into high-order structured tensors, and a kernel factor with a special structure (Hankel) is constructed to approximate and achieve tensor completion for improving the approximated accuracy. We suggested a new algorithm that effectively alleviates the burden of rank selection and improves the accuracy of tensor completion.

The rest of this paper is organized as follows: the notations used throughout this paper and related preliminaries are introduced in Section 2. Section 3 establishes the proposed algorithms in detail. Convergence analysis is discussed under mild conditions in Section 4. In Section 5, we compare the proposed algorithms with the previous algorithms to show the effectiveness of the new algorithms. Finally, we conclude the paper in Section 6.

## 2. Notation and preliminaries

Throughout this paper,  $\mathbb{N}$  stands for the set of the natural number and  $\mathbb{R}$  the real number;  $\mathbb{R}^{I_1}$  and  $\mathbb{R}^{I_1 \times I_2}$  denote the space of  $I_1$  dimensional real column vectors and  $I_1 \times I_2$  dimensional real matrices, respectively. Let  $\mathbb{R}^{I_1 \times I_2 \times \cdots \times I_N}$  be the space of Nth-order tensors with size  $I_1 \times I_2 \times \cdots \times I_N$ . The calligraphic letters, e.g.,  $\mathcal{A}$ , denote tensors; capital letters, e.g., A, denote matrices; lowercase letters, e.g., a, denote vectors; and Greek letters, e.g., a, denote scalars. For matrix X, we let  $\sigma_1(X) \geq \sigma_2(X) \geq \cdots \geq \sigma_r(X)$  be singular values of X with rank r and  $||X||_* = \operatorname{trace}((X^TX)^{\frac{1}{2}}) = \sum_{i=1}^r \sigma_i(X)$  be the nuclear norm of X, here  $\operatorname{trace}(\cdot)$  is the matrix trace operation. For an Nth-order tensor  $X \in \mathbb{R}^{I_1 \times I_2 \times \dots \times I_N}$ , its element is denoted by  $X_{i_1 i_2 \dots i_N}$ , where  $1 \leq i_n \leq I_n$  and  $1 \leq n \leq N$ . Where appropriate, a tensor sequence can also be written as [X]. The inner product of two tensors, X and Y, of the same order and size is defined as  $(X, Y) = \sum_{i_1, i_2, \dots, i_N} X_{i_1 i_2 \dots i_N} Y_{i_1 i_2 \dots i_N}$ . The Frobenius norm of a tensor is given by  $||X||_F = \langle X, X \rangle^{\frac{1}{2}}$ . The orthogonal projection operator onto an index set  $\Omega$ , denoted by  $P_{\Omega}(\cdot)$ , is defined as

$$(P_{\Omega}(X))_{i_1i_2...i_N} = \begin{cases} X_{i_1i_2...i_N}, & \text{if } i_1i_2...i_N \in \Omega, \\ 0, & \text{if } i_1i_2...i_N \in \bar{\Omega}, \end{cases} \forall X \in \mathbb{R}^{I_1 \times I_2 \times ... \times I_N},$$

where  $\bar{\Omega}$  is the complementary set of  $\Omega$ . A Hankel matrix  $H = (h_{i+j-1})_{i,j=1}^n \in \mathbb{R}^{n \times n}$  is of the form

$$H = \begin{pmatrix} h_1 & h_2 & \cdots & h_n \\ h_2 & h_3 & \cdots & h_{n+1} \\ \vdots & \vdots & \ddots & \vdots \\ h_n & h_{n+1} & \cdots & h_{2n-1} \end{pmatrix},$$

which is determined by 2n - 1 entries, that is, the first row and the last column.

The TR decomposition represents a tensor of higher order by circular multi-linear products over a sequence of low-order latent core tensors, i.e., TR factors. For n = 1, 2, ..., N, the TR factors are denoted by  $\mathcal{G}^{(n)} \in \mathbb{R}^{r_n \times I_n \times r_{n+1}}$  and each consists of two rank modes (i.e., mode-1 and mode-3) and one dimension mode (i.e., mode-2). The syntax  $r = \{r_1, r_2, ..., r_N\}$  denotes the TR-rank, which controls the model complexity of the TR decomposition. The TR decomposition applies trace operations, and all of the TR factors are set to be 3-order; thus, the TR decomposition relaxes the rank constraint on the first and last core of the TT to  $r_1 = r_{N+1}$ . Moreover, the TR decomposition linearly scales to the order of the tensor, and in this way, it overcomes the curse of dimensionality. In this case, the TR can be considered as a linear combination of the TTs and hence offers a powerful and generalized representation ability. The element-wise relation of the TR decomposition and the generated tensor is given by:

$$X_{i_1 i_2 \dots i_N} = \text{trace} \left\{ \prod_{n=1}^N G_{i_n}^{(n)} \right\},$$
 (2.1)

where  $X_{i_1i_2...i_N}$  denotes the  $(i_1,i_2,...,i_N)$ th element of the tensor,  $G_{i_n}^{(n)} \in \mathbb{R}^{r_n \times r_{n+1}}$  is the  $i_n$ -th mode-2 slice matrix of  $\mathcal{G}^{(n)}$ , which can also be denoted by  $\mathcal{G}^{(n)}(:,i_n,:)$  according to the MATLAB notation. For simplicity, we denote the TR decomposition by  $X = \Psi(\mathcal{G}^{(1)},\mathcal{G}^{(2)},...,\mathcal{G}^{(N)})$ .

There now exist three "unfold" operations unfold<sub>n</sub>( $\mathcal{T}$ ) along the n-mode of an Nth-order tensor  $\mathcal{T} \in \mathbb{R}^{I_1 \times I_2 \times \cdots \times I_N}$  are respectively defined as:  $T_{\langle n \rangle} \in \mathbb{R}^{\prod_{j=1}^n I_j \times \prod_{j=n+1}^N I_j}$ , the n-unfolding matrix (see [12, 30]) with its elements defined by  $T_{i_1 i_2 \dots i_N} = T_{\langle n \rangle}(i_1 \dots i_n, i_{n+1} \dots i_N)$ ;  $T_{(n)} \in \mathbb{R}^{I_n \times \prod_{j \neq n} I_j}$ , the classical mode-n unfolding matrix (see [8, 12, 30]) with its elements defined by  $T_{i_1 i_2 \dots i_N} = T_{(n)}(i_n, i_1 \dots i_{n-1} i_{n+1} \dots i_N)$ ;  $T_{[n]} \in \mathbb{R}^{I_n \times \prod_{j \neq n} I_j}$ , the mode-n unfolding matrix (see [12, 30]) with its elements defined by  $T_{i_1 i_2 \dots i_N} = T_{[n]}(i_n, i_{n+1} \dots i_N i_1 \dots i_{n-1})$ . The difference between the latter two types of mode-n unfolding operations lies in the ordering of indices associated with the N-1 modes, which corresponds to a specific dimensional permutation performed on  $\mathcal{T}$ . Their inverse operation "fold" is respectively defined as fold<sub>n</sub>( $T_{(n)}$ ) =  $\mathcal{T}$ , fold<sub>n</sub>( $T_{(n)}$ ) =  $\mathcal{T}$ , fold<sub>n</sub>( $T_{(n)}$ ) =  $\mathcal{T}$ . By the way,  $||\mathcal{T}||_F = ||T_{(n)}||_F$  for any  $1 \leq n \leq N$  and so on.

**Definition 1.** (see [12]) The subchain tensor by merging all cores except the *n*th core  $\mathcal{G}^{(n)}$ , i.e.,  $\mathcal{G}^{(n+1)}, \ldots, \mathcal{G}^{(N)}, \mathcal{G}^{(1)}, \ldots, \mathcal{G}^{(n-1)}$ , is denoted by  $\mathcal{G}^{\neq n} \in \mathbb{R}^{r_{n+1} \times \prod_{j \neq n} I_j \times r_n}$  whose slice matrices are defined by

$$G^{\neq n}(\overline{i_{n+1}\dots i_N i_1 i_2\dots i_{n-1}}) = \prod_{j=n+1}^N G_{i_j}^{(j)} \prod_{j=1}^{n-1} G_{i_j}^{(j)}.$$
 (2.2)

**Definition 2.** (see [33]) The *n*-mode (matrix) product of a tensor  $\mathcal{T} = (\mathcal{T}_{i_1 i_2 \cdots i_N}) \in \mathbb{R}^{I_1 \times I_2 \times \cdots \times I_N}$  with a matrix  $X = (X_{ji_n}) \in \mathbb{R}^{J \times I_n}$  is denoted by  $\mathcal{T} \times_n X$  and is of size  $I_1 \times I_2 \times \cdots \times I_{n-1} \times J \times I_{n+1} \times \cdots \times I_N$ .

Elementwise, we have

$$(\mathcal{T} \times_n X)_{i_1 \cdots i_{n-1} j i_{n+1} \cdots i_N} = \sum_{i_n=1}^{I_n} \mathcal{T}_{i_1 i_2 \cdots i_N} X_{j i_n}.$$
 (2.3)

Each mode-n fiber is multiplied by the matrix X. The idea can be expressed in terms of unfolded tensors:

$$\mathcal{Y} = \mathcal{T} \times_n X \Leftrightarrow Y_{(n)} = XT_{(n)},$$

that is to say,  $\mathcal{Y} = \text{fold}_n(Y_{(n)}) = \text{fold}_n(XT_{(n)})$ .

**Definition 3.** (see [31]) A standard delay embedding transform (VDT) of a vector  $v = (v_1, v_2, \dots, v_L)^T \in \mathbb{R}^L$  with  $\tau$  is given

$$\mathcal{H}(v) = \begin{pmatrix} v_1 & v_2 & \cdots & v_{L-\tau+1} \\ v_2 & v_3 & \cdots & v_{L-\tau+2} \\ \vdots & \vdots & \ddots & \vdots \\ v_{\tau} & v_{\tau+1} & \cdots & v_L \end{pmatrix} \in \mathbb{R}^{\tau \times (L-\tau+1)},$$

it can be considered as a Hankelization operation since  $\mathcal{H}(v)$  is a Hankel matrix.

If  $S \in \mathbb{B}^{\tau(L-\tau+1)\times L} \subset \mathbb{R}^{\tau(L-\tau+1)\times L}$  with  $\mathbb{B} = \{0,1\}$  is a Boolean or a 0-1 matrix that satisfies

$$\text{vec}(\mathcal{H}(v)) = Sv$$
,

then the standard delay embedding transform can be obtained by

$$(\mathcal{H}(v)) = \text{fold}_{(I,\tau)}(Sv),$$

where  $\operatorname{fold}_{(L,\tau)}: \mathbb{R}^{\tau(L-\tau+1)} \to \mathbb{R}^{\tau \times (L-\tau+1)}$  is a folding operator from a vector to a matrix. It should be noted that S is also a duplication matrix; the diagonal elements of  $(S^TS)$  comprise the numbers of duplications for individual elements, which are usually  $\tau \in \mathbb{N}$ , but low for marginal elements.

The inverse delay embedding transform for a Hankel matrix  $V_H$  can be given by

$$\mathcal{H}_{\tau}^{-1}(V_H) = S^{\dagger} \operatorname{vec}(V_H),$$

where  $S^{\dagger} = (S^T S)^{-1} S^T$  stands for the Moore–Penrose pseudo-inverse of S. An inverse transform of a VDT can be decomposed into duplication and folding, so the inverse transform can also be decomposed into the individual corresponding inverse transforms: a vectorization operation and the Moore–Penrose pseudo-inverse.

**Definition 4.** (see [31]) A multi-way delay-embedding transform (MDT) for an *N*th order tensor  $\mathcal{T} \in \mathbb{R}^{I_1 \times I_2 \times \cdots \times I_N}$  with  $\tau = (\tau_1, \tau_2, \dots, \tau_N) \in \mathbb{N}^N$  is defined by

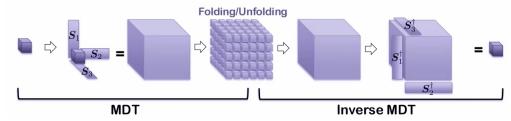
MDT: 
$$\mathcal{H}_{\tau}(\mathcal{T}) = \text{fold}_{(I,\tau)}(\mathcal{T} \times_1 S_1 \times_2 \cdots \times_N S_N),$$

where the operator  $\text{fold}_{(I,\tau)}: \mathbb{R}^{\tau_1(I_1-\tau_1+1)\times\cdots\times\tau_N(I_N-\tau_N+1)} \to \mathbb{R}^{\tau_1\times(I_1-\tau_1+1)\times\cdots\times\tau_N\times(I_N-\tau_N+1)}$  constructs a 2*N*th order tensor from the input *N*th order tensor.

The inverse MDT for a Hankel tensor is given by

invMDT: 
$$\mathcal{H}_{\tau}^{-1}(\mathcal{T}_H) = \mathcal{T}_H \times_1 S_1^{\dagger} \times \cdots \times_N S_N^{\dagger}$$

where  $\mathcal{T}_H$  is a 2Nth order tensor, this operator reshaped an Nth order tensor from the 2Nth order tensor. For convenience, they would be denoted by  $MDT(\mathcal{X})$  and  $invMDT(\mathcal{X})$  in the subsequent, respectively. In fact, the MDT is also a combination of multi-linear duplication and multi-way folding operations. Figure 1 shows flowcharts to illustrate MDT and invMDT for N = 3.



**Figure 1.** The technique of a multi-way delay-embedding transform (MDT) and its inverse for N = 3.

Let  $X \in \mathbb{R}^{I_1 \times I_2}$  be a matrix with rank r. The SVD of X is given by  $X = U\Sigma_r V^T$ , where  $U \in \mathbb{R}^{I_1 \times r}$  and  $V \in \mathbb{R}^{I_2 \times r}$  have orthogonal columns formed by the left and the right singular vectors of X, respectively, and  $\Sigma_r = \operatorname{diag}(\{\sigma_i\}_{1 \le i \le r})$  is a diagonal matrix consisting of all singular values of X, see [34]. Let  $\tau > 0$  be a parameter. The singular value shrinkage operator is given by  $D_{\tau}(X) = U\operatorname{diag}(\{\max(\sigma_i - \tau, 0)\}_{1 \le i \le r})V^T$ , see [35].

# 3. Proposed algorithm

In this section, we introduce a new algorithm by embedding a Hankel structure for low-rank tensor ring completion.

To the best of our knowledge, there are two proposed TR-based tensor completion algorithms: the TR-ALS and TR-WOPT, mentioned in above Section 1. They apply the same optimization model, which is formulated as:

$$\min_{[\mathcal{G}]} \|P_{\Omega}(\mathcal{T} - \Psi([\mathcal{G}])\|_F^2, \tag{3.1}$$

where the optimization objective is the TR factors sequence  $[\mathcal{G}]$ ,  $P_{\Omega}(\mathcal{T})$  denotes all the observed entries, the set of indices of observed entries represented by  $\Omega$ , and  $\Psi([\mathcal{G}]) = \Psi(\mathcal{G}^{(1)}, \mathcal{G}^{(2)}, \dots, \mathcal{G}^{(N)})$  denotes the approximated tensor generated by  $[\mathcal{G}]$ . Every element of  $\Psi([\mathcal{G}])$  is calculated by Eq (2.1).

Completion by rank minimization-based regularization, the model (3.1) can be formulated as:

$$\min_{X} \quad \operatorname{rank}(X) + \frac{\lambda}{2} \|P_{\Omega}(T - X)\|_{F}^{2}, \tag{3.2}$$

where X is the completed low-rank tensor, and rank(X) is a rank regularizer. The model can therefore find the low-rank structure of the data and approximate the incomplete tensor.

Because determining the tensor rank is an NP-hard problem [33, 36], work in [8, 37] extends the concept of low-rank matrix completion and defines tensor rank as a sum of the rank of the mode-*n* unfolding of the object tensor. A model of the tensor ring completion can be equivalently rewritten as:

$$\min_{\mathcal{G}} \sum_{n=1}^{N} \operatorname{rank}(\mathcal{G}^{(n)}) + \frac{\lambda}{2} \|P_{\Omega}(\mathcal{T} - \Psi([\mathcal{G}]))\|_{F}^{2},$$

Usually, the model is solved by ADMM algorithms, and it is shown to have fast convergence and good performance when the data size is small. However, when dealing with large-scale data, the multiple SVD operations in the optimization step will be intractable due to high computational cost.

By nuclear norm regularization and imposing low rankness on each of the TR factors, a tensor completion model is formulated as follows:

$$\min_{\substack{[\mathcal{G}], \mathcal{X} \\ \text{s.t.}}} \sum_{n=1}^{N} ||\mathcal{G}^{(n)}||_* + \frac{\lambda}{2} ||\mathcal{X} - \Psi([\mathcal{G}])||_F^2,$$
s.t. 
$$P_{\Omega}(\mathcal{X}) = P_{\Omega}(\mathcal{T}).$$
(3.3)

Further, the model (3.3) was finally rewritten as the tensor ring low-rank factors (TRLRF) model in [19, 33, 36, 38]:

$$\min_{\substack{[\mathcal{G}], \mathcal{X} \\ \text{s.t.}}} \sum_{n=1}^{N} \sum_{k=1}^{3} \|G_{(k)}^{(n)}\|_* + \frac{\lambda}{2} \|\mathcal{X} - \Psi([\mathcal{G}])\|_F^2, \\
\text{s.t.} \quad P_{\Omega}(\mathcal{X}) = P_{\Omega}(\mathcal{T}).$$
(3.4)

This TRLRF model has two distinctive advantages. First, the low-rank assumption is placed on tensor factors instead of on the original tensor; this greatly reduces the computational complexity of the SVD operation. Second, the low-rankness of the tensor factors can enhance the robustness to rank selection, which can alleviate the burden of searching for the optimal TR-rank and reduce the computational cost in the implementation. Immediately after, a new model is proposed that utilizes the rank relationship between a tensor and its core, which gives low-rank constraints on the two rank modes of the TR factors, i.e., the unfolding of the TR factors along mode-1 and mode-3, which can be expressed by  $\sum_{n=1}^{N} ||G_{(1)}^{(n)}||_* + \sum_{n=1}^{N} ||G_{(3)}^{(n)}||_* [20].$  Furthermore, it is noted that a slicing matrix of the mode-2 tensor core factor represents that important information of the original from definition 1, and so it can be used instead of mode-1 and mode-3 to improve computational efficiency and reduce computation time. So we developed it as follows:

$$\min_{\substack{[\mathcal{G}], \mathcal{X} \\ \text{s.t.}}} \sum_{n=1}^{N} 3 \|G_{(2)}^{(n)}\|_* + \frac{\lambda}{2} \|\mathcal{X} - \Psi([\mathcal{G}])\|_F^2, 
\text{s.t.} \quad P_{\Omega}(\mathcal{X}) = P_{\Omega}(\mathcal{T}).$$
(3.5)

By embedding the Hankel structure  $\mathcal{H}_{\mathcal{T}} = \text{MDT}(\mathcal{X})$ , the model (3.5) is reexpressed as the following model:

$$\min_{\substack{[\mathcal{G}], \mathcal{H}_{\mathcal{T}} \\ \text{s.t.}}} \sum_{n=1}^{2N} 3 \|\mathcal{G}_{(2)}^{(n)}\|_* + \frac{1}{2} \|\mathcal{H}_{\mathcal{T}} - \Psi([\mathcal{G}])\|_F^2, \\
\text{s.t.} \quad P_{\Omega}(\mathcal{X}) = P_{\Omega}(\mathcal{T}).$$
(3.6)

We apply the augmented Lagrange multiplier (ALM) method [39] to solve (3.6). By introducing  $\mathcal{M}$ ,  $\mathcal{Y}$ , then the augmented Lagrangian function optimization model of (3.6) with a constraint is given as

$$L(\mathcal{G}, \mathcal{X}, \mathcal{M}, \mathcal{Y}) = \sum_{n=1}^{2N} (3||M_{(2)}^{(n)}||_* + \langle \mathcal{Y}_n, \mathcal{M}_n - \mathcal{G}_n \rangle + \frac{\mu}{2}||\mathcal{M}_n - \mathcal{G}_n||_F^2) + \frac{\mu}{2}||\mathcal{H}_\tau - \Psi([\mathcal{G}_n])||_F^2,$$

s.t. 
$$P_{\Omega}(X) = P_{\Omega}(\mathcal{T}),$$
 (3.7)

where  $\mu > 0$  is the penalty parameter,  $\mathcal{M}_n$  is the auxiliary variable of  $\mathcal{G}_n$ ,  $\mathcal{Y}_n$  is the Lagrangian multiplier, and  $\mathcal{G}_n$ ,  $\mathcal{M}_n$ ,  $\mathcal{Y}_n$  are independent of each other. The corresponding mathematical programming can be given in the following

min 
$$L(\mathcal{G}, X, \mathcal{M}, \mathcal{Y})$$
,  
s.t.  $P_{\Omega}(X) = P_{\Omega}(\mathcal{T})$ . (3.8)

The optimization problem (3.8) is a nonconvex and nonsmooth programming. Thus, we obtain Algorithm 1.

# Algorithm 1: Embedding Structured Data in Low-rank Tensor Ring Completion.

```
Input: P_{\Omega}(\mathcal{T}), \Psi, N;
output: X
0. Let the subscripts of each element position in G_N be (i, j, k), the element values that make their
subscripts equal to sum(i + j) the element are the same. \mathcal{Y}_1 = 0, \mathcal{M}_1 = 0, \lambda = 5, \mu = 10^{-1}, \mu_{\text{max}} = 10^2,
\rho = 1.3, \epsilon = 10^{-6}, t_{\text{max}} = 5000;
1. X = P_{\Omega}(\mathcal{T})
2. \mathcal{H}_{\tau} = \text{MDT}(\mathcal{X}),
     for t = 1 : t_{\text{max}},
     while n = 1:2N
3. \mathcal{H}_{last} = \mathcal{H}_{\tau}
4. \mathcal{G}_{n+1} = \text{fold}_2 \left( (3(\mu M_{(2)}^n + Y_{(2)}^n) + \lambda H_{\text{last}} G_{<2>}^{\neq n}) (\lambda G_{<2>}^{\neq n, \text{T}} G_{<2>}^{\neq n} + 3\mu I)^{-1} \right)
5. \mathcal{M}_{n+1} = \text{fold}_2(\mathcal{D}_{\frac{1}{u}}(G_{(2)}^{n+1} - \frac{1}{u}Y_{(2)}^n))
6. \mathcal{H}_{\tau} = P_{\Omega}(\mathcal{H}_{last}) + P_{\bar{\Omega}}(\Psi([\mathcal{G}]))
7. \mathcal{Y}_{n+1} = \mathcal{Y}_n + \mu(\mathcal{M}_{n+1} - \mathcal{G}_{n+1})
8. \mu = \min(\rho \mu, \mu_{\text{max}})
9. if \frac{\|\mathcal{H}_{\tau} - \mathcal{H}_{last}\|_F}{\|\mathcal{H}_{\tau}\|_F} < \epsilon, break
     end
10. X = \text{invMDT}(\mathcal{H}_{\tau})
```

**Remark.** The algorithm 1 is conceptually quite simple, comprising three steps: (1) MDT; (2) low-rank tensor approximation; and (3) inverse MDT. In (2), we obtain a low-rank approximation of X based on the TR decomposition model (3.5).

## 4. Convergence analysis

In this section, we discuss the global convergence of Algorithm 1.

For the ease of exposition,  $\mathcal{H}$ ,  $\mathcal{M}_k$ ,  $\mathcal{G}_k$ ,  $\mathcal{Y}_k$  as described in Algorithm 1, let  $(\mathcal{M}^*, \mathcal{G}^*, \mathcal{Y}^*)$  be the solution of (3.8).

**Lemma 1.** The sequence  $\{\mathcal{Y}_k\}$  generated by Algorithm 1 is bounded.

*Proof.* Let  $\mathcal{B} = \mu_k(\mathcal{M}_{k+1} + \frac{1}{\mu_k}\mathcal{Y}_k - \mathcal{G}_k)$ . We have the following from Algorithm 1,

$$\mathcal{Y}_{k+1} = \mathcal{Y}_k + \mu_k (\mathcal{M}_{k+1} - \mathcal{G}_{k+1}) 
= \mathcal{Y}_k + \mu_k (\mathcal{M}_{k+1} - \mathcal{G}_k + \mathcal{G}_k - \mathcal{G}_{k+1}) 
= \mathcal{Y}_k + \mu_k (\mathcal{M}_{k+1} - \mathcal{G}_k) + \mu_k (\mathcal{G}_k - \mathcal{G}_{k+1}).$$

By simple deduction,

$$\mu_{k}(\mathcal{G}_{k} - \mathcal{G}_{k+1}) = \mu_{k} P_{\bar{\Omega}}(\mathcal{G}_{k} - \mathcal{G}_{k+1})$$

$$= \mu_{k} P_{\bar{\Omega}}(\mathcal{G}_{k} - (\mathcal{M}_{k+1} + \mu_{k}^{-1} \mathcal{Y}_{k}))$$

$$= P_{\bar{\Omega}}(\mu_{k}(\mathcal{M}_{k+1} + \mu_{k}^{-1} \mathcal{Y}_{k} - \mathcal{G}_{k}))$$

$$= P_{\bar{\Omega}}(\mathcal{B}).$$

And

$$\begin{split} ||\mathcal{B}||_F &= ||\mu_k(\mathcal{M}_{k+1} + \mu_k^{-1}\mathcal{Y}_k - \mathcal{G}_k)||_F \\ &= \mu_k \sum_{i=1}^N ||M_{(i)}^{k+1} + \mu_k^{-1}Y_{(i)}^k - G_{(i)}^k)||_F \\ &\leq N\mu_k ||M_{(2)}^{k+1} + \mu_k^{-1}Y_{(2)}^k - G_{(2)}^k)||_F. \end{split}$$

The truncated SVD has been used in Algorithm 1, then

$$M_{(2)}^{k+1} + \mu_k^{-1} Y_{(2)}^k = \overline{U}_k \overline{\Sigma}_k \overline{V}_k^{\mathrm{T}} + \widetilde{U}_k \widetilde{\Sigma}_k \widetilde{V}_k^{\mathrm{T}},$$

where  $\overline{U}_k \overline{\Sigma}_k \overline{V}_k^{\mathrm{T}}$  and  $\widetilde{U}_k \widetilde{\Sigma}_k \widetilde{V}_k^{\mathrm{T}}$  are both the partial SVD divided by thresholding  $\mu_k^{-1}$ , that is to say,  $\overline{\sigma}_k > \mu_k^{-1}$ ,  $\widetilde{\sigma}_k \leq \mu_k^{-1}$ .

Hence,

$$\begin{split} ||\mathcal{B}||_{F} &= ||\mu_{k}(M_{(2)}^{k+1} + \mu_{k}^{-1}Y_{(2)}^{k} - G_{(2)}^{k})||_{F} \\ &\leq N\mu_{k}||(M_{(2)}^{k+1} + \mu_{k}^{-1}Y_{(2)}^{k} - G_{(2)}^{k})||_{F} \\ &= N||\mu_{k}(\overline{U}_{k}\overline{\Sigma}_{k}\overline{V}_{k}^{T} + \widetilde{U}_{k}\widetilde{\Sigma}_{k}\widetilde{V}_{k}^{T} - U_{k}(\Sigma_{k} - \mu_{k}^{-1}I)\overline{V}_{k}^{T})||_{F} \\ &= N||\mu_{k}(\mu_{k}^{-1}\overline{U}_{k}I\overline{V}_{k}^{T} + \widetilde{U}_{k}\widetilde{\Sigma}_{k}\widetilde{V}_{k}^{T})||_{F}. \end{split}$$

From the Lemmas 1 and 4 in [39],  $\mu_k(M_{(2)}^{k+1} + \mu_k^{-1}Y_{(2)}^k - G_{(2)}^k) \in \nabla ||M_{(2)}^{k+1}||_*$ . Let  $M_{(2)}^{k+1} = U\Sigma V^{\mathrm{T}}$ ,

$$\begin{split} \|M_{(2)}^{k+1}\|_* &= \{UV^{\mathsf{T}} + W: W \in \mathbb{R}^{n \times m}, U^{\mathsf{T}}W = 0, WV = 0, \|W\|_2 \le 1\}, \\ \|UV^{\mathsf{T}} + W\|_F^2 &= \operatorname{trace}(VV^{\mathsf{T}} + W^{\mathsf{T}}W) \le N, \\ \|\mu_k(M_{(2)}^{k+1} + \mu_k^{-1}Y_{(2)}^k - G_{(2)}^k)\|_F \le \sqrt{N}, \end{split}$$

Then,

$$N||\mu_{k}(M_{(2)}^{k+1} + \mu_{k}^{-1}Y_{(2)}^{k} - G_{(2)}^{k})||_{F} \leq N\sqrt{N},$$
  
$$||\mu_{k}(\mathcal{M}_{k+1} + \mu_{k}^{-1}\mathcal{Y}_{k} - \mathcal{G}_{k})||_{F} \leq N\sqrt{N},$$
  
$$||\mathcal{B}||_{F} \leq N\sqrt{N}.$$

Thus, Lemma 1 holds.

**Theorem 1.** Suppose that the series  $\sum_{k=1}^{+\infty} \mu_k^{-1}$  diverges if  $\mu_k \to +\infty$ . Then the sequences  $\{\mathcal{M}_k\}, \{\mathcal{G}_k\}$  generated by Algorithm 1 converge to the optimal solution of (3.8).

Proof. According to Algorithm 1, we obtain

$$\mathcal{Y}_{k+1} = \mathcal{Y}_k + \mu_k (\mathcal{M}_{k+1} - \mathcal{G}_{k+1}),$$

$$\mu_k^{-1} (\mathcal{Y}_{k+1} - \mathcal{Y}_k) = \mathcal{M}_{k+1} - \mathcal{G}_{k+1},$$

$$\lim_{k \to +\infty} \mu_k^{-1} (\mathcal{Y}_{k+1} - \mathcal{Y}_k) = \lim_{k \to +\infty} (\mathcal{M}_{k+1} - \mathcal{G}_{k+1}) = 0.$$

Let  $\mathcal{M}^*$ ,  $\mathcal{G}^*$  denote the optimal solution of (3.8); this means  $\mathcal{M}^* = \mathcal{G}^*$ . In order to prove this theorem, we now consider the formula

$$\begin{split} \|\mathcal{M}_{k+1} - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}^*\|_F^2 \\ &= \|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2 - \|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 \\ &- \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2 - 2\mu_k^{-1} \langle \mathcal{G}_{k+1} - \mathcal{G}^*, \widehat{\mathcal{Y}}_{k+1} - \mathcal{Y}^* \rangle. \end{split}$$

Suppose that  $\widehat{\mathcal{Y}}_{k+1} = \mathcal{Y}_k + \mu_k(\mathcal{M}_{k+1} - \mathcal{G}_k), \mathcal{Y}^*$  is the optimal solution of the dual problem in [39].

$$\begin{split} \|\mathcal{M}_{k+1} - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}^*\|_F^2 \\ &= \|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2 - (\|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2) \\ &+ 2\langle \mathcal{M}_{k+1} - \mathcal{M}^*, \mathcal{M}_{k+1} - \mathcal{M}_k \rangle + \mu_k^{-2} \langle \mathcal{Y}_{k+1} - \mathcal{Y}^*, \mathcal{Y}_{k+1} - \mathcal{Y}_k \rangle \\ &= \|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2 - (\|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2) \\ &+ 2\langle \mathcal{M}_{k+1} - \mathcal{M}^*, \mathcal{M}_{k+1} - \mathcal{M}_k \rangle + 2\mu_k^{-1} \langle \mathcal{Y}_{k+1} - \mathcal{Y}^*, \mathcal{M}_{k+1} - \mathcal{G}_{k+1} \rangle \\ &= \|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2 - (\|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2) \\ &+ 2\langle \mathcal{M}_{k+1} - \mathcal{M}^*, \mathcal{M}_{k+1} - \mathcal{M}_k \rangle + 2\mu_k^{-1} (\langle \mathcal{Y}_{k+1} - \mathcal{Y}^*, \mathcal{M}_{k+1} - \mathcal{M}^* \rangle \\ &+ \langle \mathcal{Y}_{k+1} - \mathcal{Y}^*, \mathcal{G}^* - \mathcal{G}_{k+1} \rangle) \\ &= \|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2 - (\|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2) \\ &- 2\mu_k^{-1} \langle \mathcal{G}_{k+1} - \mathcal{G}^*, \widehat{\mathcal{Y}}_{k+1} - \mathcal{Y}^* \rangle. \end{split}$$

Since  $||\mathcal{G}||_*$  is a convex function and  $\widehat{\mathcal{Y}}_{k+1} \in \nabla ||\mathcal{G}_{k+1}||_*$ , then

$$\langle \mathcal{G}_{k+1} - \mathcal{G}^*, \widehat{\mathcal{Y}}_{k+1} - \mathcal{Y}^* \rangle \geq 0.$$

It is noted that  $\|\mathcal{M}_{k+1} - \mathcal{M}_k\|_F^2 + \mu_k^{-2} \|\mathcal{Y}_{k+1} - \mathcal{Y}_k\|_F^2 \ge 0$ , and

$$||\mathcal{M}_{k+1} - \mathcal{M}^*||_F^2 + \mu_k^{-2}||\mathcal{Y}_{k+1} - \mathcal{Y}^*||_F^2 \le ||\mathcal{M}_k - \mathcal{M}^*||_F^2 + \mu_k^{-2}||\mathcal{Y}_k - \mathcal{Y}^*||_F^2 + 2\mu_k^{-1}\langle \mathcal{G}_{k+1} - \mathcal{G}^*, \widehat{\mathcal{Y}}_{k+1} - \mathcal{Y}^* \rangle.$$

Then  $\{\|\mathcal{M}_k - \mathcal{M}^*\|_F^2 + \mu_k^{-2}\|\mathcal{Y}_k - \mathcal{Y}^*\|_F^2\}$  is not increasing, and  $\sum_{k=1}^{\infty} \mu_k^{-1} \langle \mathcal{M}_k - \mathcal{M}^*, \mathcal{Y}_k - \mathcal{Y}^* \rangle < +\infty$ . Consequently,  $\mathcal{M}^*, \mathcal{G}^*$  is the optimal solution of (3.7) by the proof of Lemma 2 in [39]. So the theorem is proven.

**Theorem 2.** Assume that the tensor  $\mathcal{H}_{\tau}$  is formed by  $[\mathcal{G}_k]$ . Then

$$\|\mathcal{H}_{\tau} - \mathcal{H}^*\|_F \leq \|\Psi([\mathcal{G}_k]) - \mathcal{H}^*\|_F$$

here  $\mathcal{H}^*$  is the solution of (3.6). *Proof.* Because of

$$\begin{split} \|\Psi([\mathcal{G}_{k}]) - \mathcal{H}^{*}\|_{F}^{2} &= \|\Psi([\mathcal{G}_{k}]) - \mathcal{H}_{\tau} + \mathcal{H}_{\tau} - \mathcal{H}^{*}\|_{F}^{2} \\ &= \|\Psi([\mathcal{G}_{k}]) - \mathcal{H}_{\tau}\|_{F}^{2} + \|\mathcal{H}_{\tau} - \mathcal{H}^{*}\|_{F}^{2} + 2\langle\Psi([\mathcal{G}_{k}]) - \mathcal{H}_{\tau}, \mathcal{H}_{\tau} - \mathcal{H}^{*}\rangle \\ &= \|\Psi([\mathcal{G}_{k}]) - \mathcal{H}_{\tau}\|_{F}^{2} + \|\mathcal{H}_{\tau} - \mathcal{H}^{*}\|_{F}^{2}, \end{split}$$

then  $\|\mathcal{H}_{\tau} - \mathcal{H}^*\|_F \le \|\Psi([\mathcal{G}_k]) - \mathcal{H}^*\|_F$  holds.

The theorem is proven.

## 5. Numerical experiments

In this section, we conducted experiments to testify to the effectiveness of our algorithm 1 by comparing some algorithms in the relative square error (RSE), peak signal-to-noise ratio (PSNR), the computing time in seconds (CPU(s)), and structural similarity (SSIM) [40]. They are defined as

$$PSNR = 10 \log_{10} (255^2 / MSE),$$

where the mean squared error (MSE) was defined by

$$MSE = \frac{\|X - \mathcal{T}\|_F^2}{\text{num}(\mathcal{T})}, \quad \text{num}(\mathcal{T}) \text{ means the number of elements of } \mathcal{T};$$

$$RSE = \frac{\|\mathcal{X} - \mathcal{T}\|_F}{\|\mathcal{T}\|_F};$$

$$SSIM(x, y) = [(l(x, y)]^{\alpha} \cdot [(c(x, y)]^{\beta} \cdot [(s(x, y)]^{\gamma},$$

where  $\alpha, \beta, \gamma$  are all positive real numbers. As for the details of the functionals l(x, y), c(x, y), and s(x, y), please refer to [40].

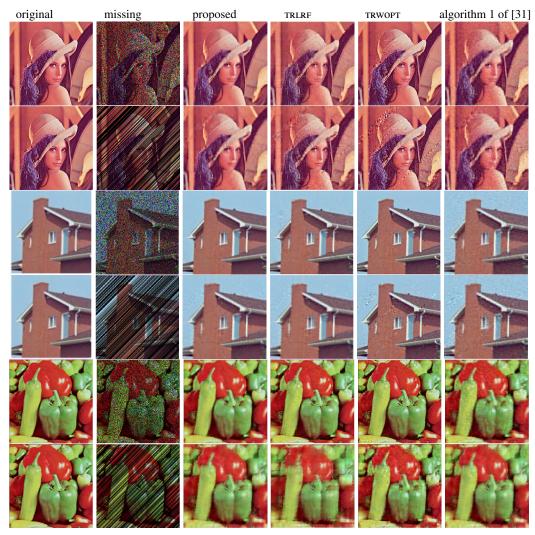
The TR-WOPT [20] and TR-LRF [38]) algorithms were used to solve the model (3.4); algorithm 1 from [31] was used to solve the model (20) in [31]; and our algorithm 1 was exploited to solve the model (3.6). In the experiments, we selected three benchmark images, named Lena, House, and Peppers with the size of them being  $256 \times 256 \times 3$ , which can be considered a 3rd order tensor. p represents the sampling rate in the test.

There are three cases for the experiment in the following. In our implementations, all the codes were written by Matlab R2019b and run on a PC with an Intel Xeon E5 processor, 2.5GHz, and 20GB memory.

# 5.1. Verification of the proposed algorithm using benchmark images inpainting

We prepared two missing formats of the images. One of them is randomly missing, and the other is missing with special structures. The sampling rate of those images was set as p = 0.5, a light blurring. Figure 2 and Table 1 have reported the experimental results obtained after the completion of three incomplete images for four algorithms. Evidently, the result obtained by the proposed algorithm 1 was better than that by other algorithms. Table 1 shows the SSIM, PSNR, RSE, and CPU(s) for these comparisons, where the best RSE, PSNR, CPU(s) and SSIM values are emphasized in bold font.

Furthermore, it is noted that all four algorithms have completed the filling task well for two missing formats by Figure 2 and Table 1. The proposed algorithm and algorithm 1 of [31] exhibit better visual effects for the structured missing case *facade 2*. Due to their inclusion of MDT, their advantage lies not in computation time but in the completion effect.



**Figure 2.** Color images completed with various algorithms. Two missing color images were artificially generated comprising: random missing case *facade 1*, structured missing case *facade 2*.

<b>Table 1.</b> Computational results for the missing images.							
image	missing	item	proposed	TRLRF	TRWOPT	alg.1 of [31]	
		SSIM	0.8790	0.7850	0.7650	0.8025	
	facade 1	PSNR	27.028	25.373	26.217	26.572	
		RSE	0.0800	0.0970	0.0920	0.0730	
Lena		CPU(s)	8.4800	6.3297	6.4905	8.4411	
		SSIM	0.8530	0.7470	0.6130	0.7344	
	facade 2	PSNR	25.665	24.495	22.181	24.027	
		RSE	0.0920	0.1080	0.1470	0.1338	
		CPU(s)	6.1768	6.97 50	6.8912	6.1020	
		SSIM	0.9030	0.8730	0.8170	0.8820	
	facade 1	PSNR	27.587	26.922	26.913	27.307	
		RSE	0.0710	0.0750	0.0770	0.0723	
House		CPU(s)	9.0821	6.4911	6.5720	9.1208	
		SSIM	0.8580	0.8030	0.6870	0.8117	
	facade 2	PSNR	26.937	25.451	24.313	26.229	
		RSE	0.0790	0.0890	0.1050	0.0823	
		CPU(s)	6.3728	7.0957	7.1120	6.4432	
		SSIM	0.8730	0.8010	0.7130	0.8331	
	facade 1	PSNR	26.678	25.449	24.899	26.187	
		RSE	0.0920	0.1070	0.1260	0.0923	
Peppers		CPU(s)	11.0748	9.2470	9.4965	10.3912	
		SSIM	0.8480	0.7560	0.5920	0.8510	
	facade 2	PSNR	26.321	24.285	21.691	26.556	
		RSE	0.1070	0.1360	0.1760	0.1121	
		CPU(s)	10.3477	9.0231	9.0020	10.0015	

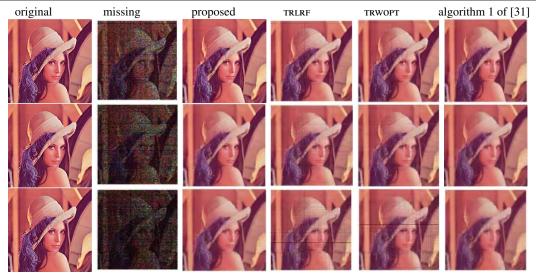
# 5.2. Comparison using benchmark images inpainting for row/column fibres missing

In order to compare the performance of the new algorithm in solving images inpainting with row/column fibers missing, we prepared three row/column fibers missing formats of the images mentioned above for this experiment, which are shown in Figures 3-5. The missing rates of the image were set as p = 0.3, 0.4, 0.5, and the corresponding missing fibers in the row and column are  $p_r = 20, 10, 5$  and  $p_c = 24, 12, 6$ , respectively.

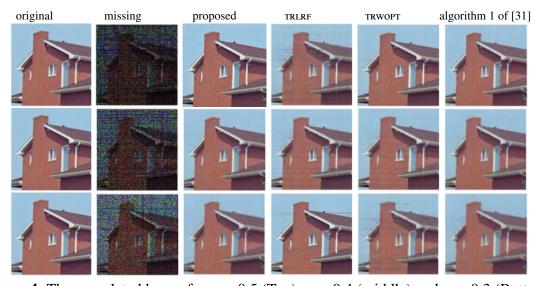
From the visual effect of Figures 3–5 and the results recorded in Table 2. It is clear that the result obtained by our algorithm 1 and the algorithm 1 of [31] were much better than that of the other two algorithms without MDT. That is to say, our algorithm 1 and the algorithm 1 of [31] are more suitable for special cases of the tensor completion. Table 2 shows the SSIM, PSNR, and RSE for these comparisons, where the best RSE, PSNR, and SSIM values are emphasized in bold font.

**Table 2.** Completed results of three images for the different missing rate.

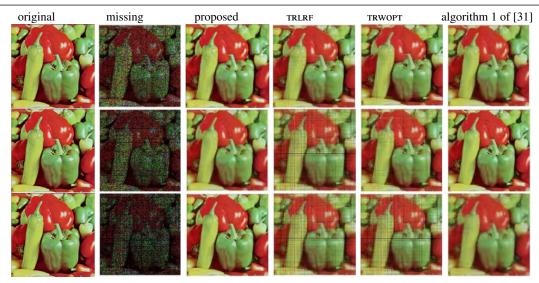
						8
image	missing	item	proposed	TRLRF	TRWOPT	alg.1 of [31]
Lena	p=0.3	SSIM	0.9499	0.9303	0.9039	0.9019
	$p_r = 20$	<b>PSNR</b>	24.2172	23.6465	22.3610	22.2662
	$p_c = 24$	RSE	0.0920	0.1388	0.1403	0.1196
	p=0.4	SSIM	0.9630	0.9394	0.9383	0.9525
	$p_r = 10$	<b>PSNR</b>	26.181	25.1286	25.0333	25.7863
	$p_c = 12$	RSE	0.0870	0.1014	0.1024	0.0936
	p=0.5	SSIM	0.9733	0.8717	0.8130	0.9721
	$p_r$ =5	<b>PSNR</b>	28.7825	28.1753	28.0181	28.7815
	$p_c$ =6	RSE	<b>0.0606</b> 8	0.0693	0.0740	0.0636
	p=0.3	SSIM	0.8808	0.8538	0.8525	0.8170
House	$p_r = 20$	<b>PSNR</b>	24.7000	23.5651	23.5148	24.0913
	$p_c = 24$	RSE	0.0963	0.1180	0.1186	0.1007
	p=0.4	SSIM	0.9179	0.8820	0.8819	0.9670
	$p_r = 10$	<b>PSNR</b>	25.3162	25.0296	25.0206	25.4313
	$p_c = 12$	RSE	0.0809	0.0923	0.0894	0.0795
	p=0.5	SSIM	0.9583	0.9417	0.9416	0.9430
	$p_r$ =5	<b>PSNR</b>	27.9485	27.6135	27.6093	27.7811
	$p_c$ =6	RSE	0.0521	0.0683	0.0683	0.0570
Peppers	p=0.3	SSIM	0.9818	0.8843	0.8847	0.9130
	$p_r = 20$	<b>PSNR</b>	30.3795	22.3312	22.3587	28.8990
	$p_c$ =24	RSE	0.0551	0.1393	0.1388	0.0860
	p=0.4	SSIM	0.9218	0.8824	0.9188	0.9208
	$p_r=10$	<b>PSNR</b>	24.5184	22.0556	23.4397	24.691
	$p_c = 12$	RSE	0.1124	0.1576	0.1393	0.1060
	p=0.5	SSIM	0.9588	0.9237	0.9528	0.9536
	$p_r$ =5	<b>PSNR</b>	25.9665	24.8717	25.1861	25.7920
	$p_c$ =6	RSE	0.0818	0.1104	0.1114	0.0847
						<del></del>



**Figure 3.** The completed Lena for p = 0.5 (Top), p = 0.4 (middle) and p = 0.3 (Bottom).



**Figure 4.** The completed house for p = 0.5 (Top), p = 0.4 (middle) and p = 0.3 (Bottom).

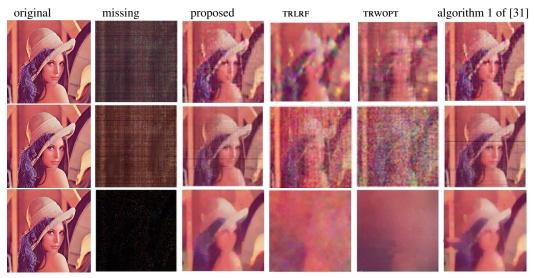


**Figure 5.** The completed peppers for p = 0.5 (Top), p = 0.4 (middle) and p = 0.3 (Bottom).

## 5.3. Comparison using Lena's image for missing slice

To compare the performance of the proposed algorithm with the other algorithms for solving image completion problems with a slice missing, we prepared Lena's images with badly random missing; the lower sampling rates of the image were badly set as 0.01, 0.03, and 0.05, respectively.

From the visual effect of Figure 6 and the results recorded in Table 3. It is noted that the results obtained by our algorithm 1 and algorithm 1 of [31] were much better than those by two other algorithms without MDT. That is to say, our algorithm 1 and algorithm 1 of [31] are more suitable for this special case of the tensor completion problems. Table 3 shows the SSIM, PSNR, and RSE for these comparisons, where the best RSE, PSNR, and SSIM values are emphasized in bold font.



**Figure 6.** The completed Lena for p = 0.05 (Top), p = 0.03 (middle) and p = 0.01 (Bottom).

missing	item	proposed	TRLRF	TRWOPT	alg.1 of [31]
p=0.05	SSIM	0.6729	0.5720	0.5581	0.6631
	<b>PSNR</b>	21.5228	18.0491	19.6552	21.0217
	RSE	0.5080	0.7901	0.8312	0.5126
p=0.03	SSIM	0.8530	0.6470	0.5840	0.8130
	<b>PSNR</b>	22.6065	19.4995	18.9337	21.9181
	RSE	0.0920	0.8080	0.8869	0.1014
p=0.01	SSIM	0.8530	0.4470	0.4584	0.8555
	<b>PSNR</b>	16.665	14.4956	13.9897	16.181
	RSE	0.2920	0.9780	0.9921	0.3470

**Table 3.** Completed results of missing Lena for the different missing rate.

## 6. Conclusions

This paper focuses on the problem of completing a tensor with structured missing or missing slices or missing row/column fibers. In the proposed algorithm, embedding structured data in a tensor for solving a non-convex optimization model of tensor completion problems by multi-way delay-embedding transform (MDT) makes the tensor to be completed with a special structure. We have introduced an efficient and high-performance tensor completion algorithm based on the tensor ring rank, constructing latent tensor ring factors with Hankel structure the approximates the original tensor starting from the tensor structure. It has been shown to effectively handle model selection and provide much better completion results. This has been demonstrated from the extensive experiments on images inpainting.

The proposed algorithm has many potential extensions, such as using different embedding transformations in step 1 of our algorithm 1 or different approximated methods in step 2 of our algorithm 1.

## **Authors' contributions**

Ruiping Wen: Conceptualization, Writing-original draft, Writing-review & editing; Tingyan Liu: Writing-original draft, Writing-review & editing; Yalei Pei: Writing-original draft, Writing-review & editing. All authors read and approved the final manuscript.

#### **Use of Generative-AI tools declaration**

The authors declare they have not used Artificial Intelligence (AI) tools in the creation of this article.

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## **Conflict of interest**

The authors declare that they have no conflict of interests.

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