

# Algorithms for Nonnegative Tensor Factorization

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# Algorithms for Nonnegative Tensor Factorization

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## ABSTRACT

Nonnegative Matrix Factorization (NMF) is an efficient technique to approximate a large matrix containing only nonnegative elements as a product of two nonnegative matrices of significantly smaller size. The guaranteed nonnegativity of the factors is a distinctive property that other widely used matrix factorization methods do not have.

Matrices can also be seen as second-order tensors. For some problems, it is necessary to process tensors of third or higher order. For this purpose, NMF can be generalized to Nonnegative Tensor Factorization (NTF). NMF and NTF are used in various application areas, for example in document classification and multi-way data analysis.

The aim of this report is to give an overview over some algorithms to compute Nonnegative Tensor Factorizations, including two multiplicative algorithm based on the Alpha-divergence and the Beta-divergence, respectively, two Hierarchical Alternating Least Squares algorithms and a Block Principal Pivoting algorithm utilizing matricization.

## KEY WORDS

High Performance Computing, Nonnegative Tensor Factorization, Nonnegative Matrix Factorization

## 1 Introduction

A characteristic property of modern society is the increasing need to process large amounts of data. One important class of data is represented by nonnegative matrices and tensors, which occur in many application areas. These are often considerably large, which makes their processing and evaluation difficult and time-consuming.

Nonnegative Matrix Factorization, (abbreviated as NMF or NNMF), is a technique to approximate a nonnegative matrix as a product of two nonnegative matrices. The two resulting matrices are usually smaller than the original matrix and therefore easier to handle and process. In the last decade, NMF has become quite popular and has been applied to a wide variety of practical problems.

The idea of such a factorization was published in 1994 under the name “Positive Matrix Factorization” [21]. In 1999, an article in Nature [14] about Nonnegative Matrix Factorization caught the attention of a wide audience. Several papers were written about NMF since then, discussing its properties, algorithms, modifications and often also possible applications. Some of the various areas where Nonnegative Matrix Factorization was successfully applied are text mining [23] [24] [1], classification of documents [2] and emails [10], clustering [28] [18], spectral data analysis [11] [22] [1], face recognition [29], distance estimation in networks [19], the analysis of EEG data [16], separation of sound sources [27], music transcription [25] [26], computational biology, for example molecular pattern discovery and class comparison and prediction [3] [8] [7] and neuroscience [6].

In contrast to other methods such as singular value decomposition (SVD) or principal component analysis (PCA), NMF has the distinguishing property that the factors are guaranteed to be nonnegative, which allows to view the factorization as an additive combination of features.

## 2 The NMF Problem

An informal description of the NMF problem is: Given a nonnegative matrix  $\mathbf{Y}$  of size  $I \times T$ , find two nonnegative matrices  $\mathbf{A}$  (size  $I \times J$ ) and  $\mathbf{X}$  (size  $J \times T$ ) such that their product  $\mathbf{AX}$  approximates  $\mathbf{Y}$ . Figure 1 illustrates the NMF problem for  $I = 7$ ,  $T = 9$  and  $J = 3$ .

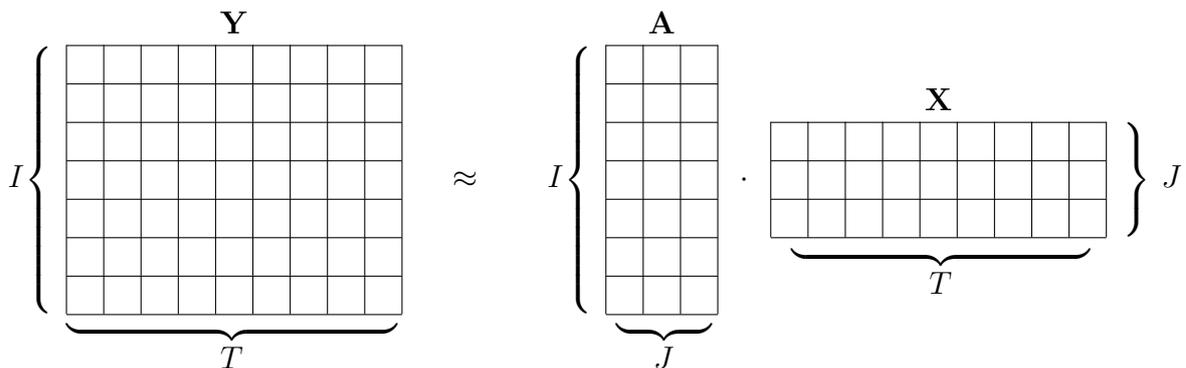


Figure 1: Illustration of the NMF problem

A matrix is called nonnegative if all its elements are  $\geq 0$ . In practical cases, the chosen  $J$  is usually much smaller than  $I$  and  $T$ . It should be noted that, in general, it is not

possible to find  $\mathbf{A}$  and  $\mathbf{X}$  such that  $\mathbf{AX} = \mathbf{Y}$ . Hence, NMF is “only” an approximation, for this reason it is sometimes called Approximative Nonnegative Matrix Factorization or Nonnegative Matrix Approximation. This is sometimes expressed as  $\mathbf{Y} = \mathbf{AX} + \mathbf{E}$ , where  $\mathbf{E}$  is a matrix of size  $I \times T$  that represents the approximation error. Thus,  $\mathbf{AX}$  can be seen as a compressed representation of  $\mathbf{Y}$ , with a rank of  $J$  or less.

Formally, NMF can be defined as [20]:

**Definition (NMF).** *Given a nonnegative matrix  $\mathbf{Y} \in \mathbb{R}^{I \times T}$  and a positive integer  $J$ , find nonnegative matrices  $\mathbf{A} \in \mathbb{R}^{I \times J}$  and  $\mathbf{X} \in \mathbb{R}^{J \times T}$  that minimize the functional*

$$f(\mathbf{A}, \mathbf{X}) = \frac{1}{2} \|\mathbf{Y} - \mathbf{AX}\|_F^2.$$

In [20],  $J < \min\{I, T\}$  is explicitly required, this is not strictly necessary, but true in almost all practical cases. For an  $I \times T$  matrix  $\mathbf{M}$ ,  $\|\mathbf{M}\|_F$  is the Frobenius norm of  $\mathbf{M}$ , defined as

$$\|\mathbf{M}\|_F := \sqrt{\sum_{i=1}^I \sum_{t=1}^T m_{i,t}^2}$$

where  $m_{i,t}$  denotes the element of  $\mathbf{M}$  with row index  $i$  and column index  $t$ . Therefore,  $f(\mathbf{A}, \mathbf{X})$  is the square of the Euclidean distance between  $\mathbf{Y}$  and  $\mathbf{AX}$  with an additional factor  $\frac{1}{2}$ . The problem is convex in  $\mathbf{A}$  and in  $\mathbf{X}$  separately, but not in both simultaneously [9].

We note that it is also possible to use other measures to express the distance between  $\mathbf{Y}$  and  $\mathbf{AX}$ , for example the Kullback-Leibler divergence [15], Csiszár’s divergences [5] or Alpha- and Beta-divergences [6]. Different measures yield different NMF algorithms, or at least different update steps for the algorithms.

Every column of  $\mathbf{A}$  can be interpreted as a basis feature of size  $I$ . In total,  $\mathbf{A}$  contains  $J$  basis features. The multiplication of  $\mathbf{A}$  with the nonnegative matrix  $\mathbf{X}$  yields a nonnegative matrix  $\mathbf{AX}$ , where every column of  $\mathbf{AX}$  is an additive (or non-subtractive) combination of weighted basis features (columns of  $\mathbf{A}$ ). The famous paper on NMF in Nature [14] uses NMF to represent faces as additive combinations of local parts such as eyes, nose, mouth, etc. However, it was shown in [17] that NMF does not always find such localized features.

## 3 The NTF Problem

Matrices are second-order tensors. For some applications, for example in multi-way data analysis, the input data are tensors of third or higher order. Therefore, it is desirable to generalize Nonnegative Matrix Factorization to Nonnegative Tensor Factorization.

### 3.1 Notation

For the formulation of the NTF problem and the algorithms, the following symbols are used:

- $\circ$  outer product
- $\odot$  Khatri-Rao product
- $\otimes$  Hadamard product
- $\oslash$  element-wise division
- $\times_n$  mode- $n$  product of tensor and matrix
- $\mathbf{A}^{\odot-n} = \mathbf{A}^{(N)} \odot \dots \odot \mathbf{A}^{(n+1)} \odot \mathbf{A}^{(n-1)} \odot \dots \odot \mathbf{A}^{(1)}$

### 3.2 Problem definition

The Nonnegative Tensor Factorization problem can be formulated as nonnegative canonical decomposition / parallel factor decomposition (CANDECOMP / PARAFAC) as follows (after [6]):

**Definition (NTF).** *Given an  $N$ -th order tensor  $\underline{\mathbf{Y}} \in \mathbb{R}^{I_1 \times I_2 \times \dots \times I_N}$  and a positive integer  $J$ , factorize  $\underline{\mathbf{Y}}$  into a set of  $N$  nonnegative component matrices  $\mathbf{A}^{(n)} = [\mathbf{a}_1^{(n)}, \mathbf{a}_2^{(n)}, \dots, \mathbf{a}_J^{(n)}] \in \mathbb{R}^{I_n \times J}$ , ( $n = 1, 2, \dots, N$ ) representing the common (loading) factors, that is,*

$$\underline{\mathbf{Y}} = \hat{\underline{\mathbf{Y}}} + \underline{\mathbf{E}} = \sum_{j=1}^J \mathbf{a}_j^{(1)} \circ \mathbf{a}_j^{(2)} \circ \dots \circ \mathbf{a}_j^{(N)} + \underline{\mathbf{E}} =$$

$$\mathbf{I} \times_1 \mathbf{A}^{(1)} \times_2 \mathbf{A}^{(2)} \dots \times_N \mathbf{A}^{(N)} + \underline{\mathbf{E}} = \llbracket \mathbf{A}^{(1)}, \mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)} \rrbracket + \underline{\mathbf{E}}$$

with  $\|\mathbf{a}_j^{(n)}\|_2 = 1$  for  $n = 1, 2, \dots, N-1$  and  $j = 1, 2, \dots, J$ .

The tensor  $\underline{\mathbf{E}}$  is the approximation error. Figure 2 illustrates the decomposition for a third-order tensor.

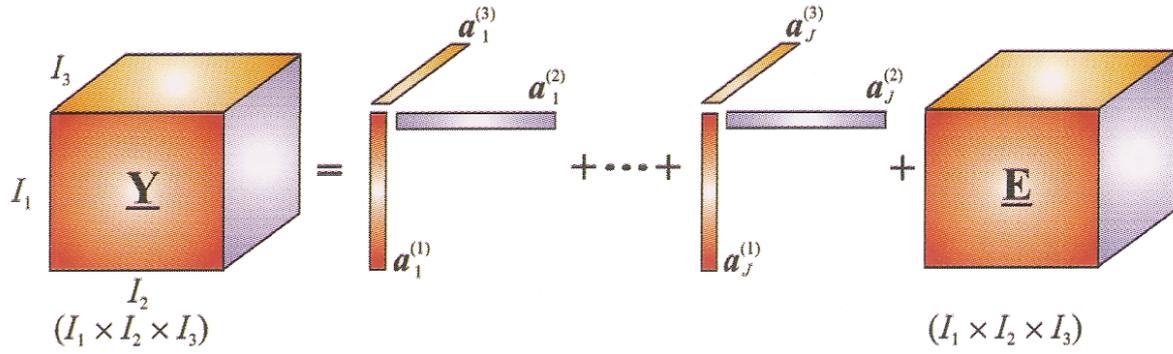


Figure 2: The NTF model for a third-order tensor (from [6])

## 4 Alpha NTF

The first algorithm presented here uses multiplicative updates based on the Alpha-divergence [6]. Multiplicative algorithms are relatively simple, but can be slower in comparison to enhanced HALS algorithms.

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**Algorithm 1:** Alpha NTF (from [6])

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**Input:**  $\underline{\mathbf{Y}}$ : input data of size  $I_1 \times I_2 \times \dots \times I_N$ ,  $J$ : number of basis components

**Output:**  $N$  component matrices  $\mathbf{A}^{(n)} \in \mathbb{R}_+^{I_n \times J}$

```

1 begin
2   ALS or random initialization for all factors  $\mathbf{A}^{(n)}$ ;
3    $\mathbf{A}_l^{(n)} = \mathbf{A}^{(n)} \text{diag}\{\mathbf{1}^T \mathbf{A}^{(n)}\}^{-1}$  for  $\forall n$ ;      /* normalize to unit length */
4    $\mathbf{A}^{(n)} = \mathbf{A}_l^{(n)}$  for  $\forall n \neq N$ ;
5   repeat
6      $\hat{\underline{\mathbf{Y}}} = \llbracket \mathbf{A}^{(1)}, \mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)} \rrbracket$ ;
7     for  $n = 1$  to  $N$  do
8        $\mathbf{A}^{(n)} \leftarrow \mathbf{A}^{(n)} \otimes \left[ \left( \underline{\mathbf{Y}}_{(n)} \oslash \hat{\underline{\mathbf{Y}}}_{(n)} \right)^{[\alpha]} \mathbf{A}_l^{\odot -n} \right]^{[1/\alpha]}$ ;
9        $\mathbf{A}_l^{(n)} = \mathbf{A}^{(n)} \text{diag}\{\mathbf{1}^T \mathbf{A}^{(n)}\}^{-1}$ ;      /* normalize to unit length */
10      if  $n \neq N$  then
11         $\mathbf{A}^{(n)} = \mathbf{A}_l^{(n)}$ ;
12      end
13    end
14  until a stopping criterion is met;
15 end
```

---

## 5 Beta NTF

It is also possible to formulate a multiplicative Nonnegative Tensor Factorization algorithm based on the Beta-divergence [6]:

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**Algorithm 2:** Beta NTF (from [6])

---

**Input:**  $\underline{\mathbf{Y}}$ : input data of size  $I_1 \times I_2 \times \dots \times I_N$ ,  $J$ : number of basis components  
**Output:**  $N$  component matrices  $\mathbf{A}^{(n)} \in \mathbb{R}_+^{I_n \times J}$

```

1 begin
2   ALS or random initialization for all factors  $\mathbf{A}^{(n)}$ ;
3   repeat
4      $\hat{\underline{\mathbf{Y}}} = \llbracket \mathbf{A}^{(1)}, \mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)} \rrbracket$ ;
5     for  $n = 1$  to  $N$  do
6        $\mathbf{A}^{(n)} \leftarrow \mathbf{A}^{(n)} \circledast \left[ \left( \mathbf{Y}_{(n)} \circledcirc \hat{\mathbf{Y}}_{(n)}^{[\beta-1]} \right) \mathbf{A}^{\odot-n} \right] \circledcirc \left( \hat{\mathbf{Y}}_{(n)}^{[\beta]} \mathbf{A}^{\odot-n} \right)$ ;
7       if  $n \neq N$  then
8          $\mathbf{A}^{(n)} \leftarrow \mathbf{A}^{(n)} \text{diag}\{\mathbf{1}^T \mathbf{A}^{(n)}\}^{-1}$ ;           /* normalize */
9       end
10    end
11  until a stopping criterion is met;
12 end
```

---

## 6 HALS NTF

HALS stands for Hierarchical Alternating Least Squares. The idea is to minimize a set of local cost functions with the same global minima by approximating rank-one tensors [4]. Two algorithms are presented here, first a simple HALS algorithm (Algorithm 3) and then another algorithm that reduces computation of the expensive Khatri-Rao products (Algorithm 4) [6]. Both algorithms use the squared Euclidean distance, but it is also possible to formulate HALS algorithms based on Alpha- and Beta-divergences.

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**Algorithm 3:** Simple HALS NTF (from [6])

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**Input:**  $\underline{\mathbf{Y}}$ : input data of size  $I_1 \times I_2 \times \dots \times I_N$ ,  $J$ : number of basis components

**Output:**  $N$  component matrices  $\mathbf{A}^{(n)} \in \mathbb{R}_+^{I_n \times J}$

```
1 begin
2   ALS or random initialization for all factors  $\mathbf{A}^{(n)}$ ;
3    $\mathbf{a}_j^{(n)} \leftarrow \mathbf{a}_j^{(n)} / \|\mathbf{a}_j^{(n)}\|_2$  for  $\forall j, n = 1, 2, \dots, N - 1$ ;      /* normalize to unit
   length */
4    $\underline{\mathbf{E}} = \underline{\mathbf{Y}} - \hat{\underline{\mathbf{Y}}} = \underline{\mathbf{Y}} - \llbracket \{\mathbf{A}\} \rrbracket$ ;      /* residual tensor */
5   repeat
6     for  $j = 1$  to  $J$  do
7        $\underline{\mathbf{Y}}^{(j)} = \underline{\mathbf{E}} + \llbracket \mathbf{a}_j^{(1)}, \mathbf{a}_j^{(2)}, \dots, \mathbf{a}_j^{(N)} \rrbracket$ ;
8       for  $n = 1$  to  $N$  do
9          $\mathbf{a}_j^{(n)} \leftarrow \left[ \mathbf{Y}_{(n)}^{(j)} \{ \mathbf{a}_j^{\odot -n} \} \right]_+$ ;
10        if  $n \neq N$  then
11           $\mathbf{a}_j^{(n)} \leftarrow \mathbf{a}_j^{(n)} / \|\mathbf{a}_j^{(n)}\|_2$ ;      /* normalize to unit length */
12        end
13      end
14       $\underline{\mathbf{E}} = \underline{\mathbf{Y}}^{(j)} - \llbracket \mathbf{a}_j^{(1)}, \mathbf{a}_j^{(2)}, \dots, \mathbf{a}_j^{(N)} \rrbracket$ ;
15    end
16  until a stopping criterion is met;
17 end
```

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---

**Algorithm 4:** FAST HALS NTF (from [6])

---

**Input:**  $\mathbf{Y}$ : input data of size  $I_1 \times I_2 \times \dots \times I_N$ ,  $J$ : number of basis components

**Output:**  $N$  component matrices  $\mathbf{A}^{(n)} \in \mathbb{R}_+^{I_n \times J}$

```
1 begin
2   ALS or random initialization for all factors  $\mathbf{A}^{(n)}$ ;
3    $\mathbf{a}_j^{(n)} \leftarrow \mathbf{a}_j^{(n)} / \|\mathbf{a}_j^{(n)}\|_2$  for  $\forall j, n = 1, 2, \dots, N - 1$ ;      /* normalize to unit
   length */
4    $\mathbf{T}^{(1)} = (\mathbf{A}^{(1)T} \mathbf{A}^{(1)}) \otimes \dots \otimes (\mathbf{A}^{(N)T} \mathbf{A}^{(N)})$ ;
5   repeat
6      $\gamma = \text{diag}(\mathbf{A}^{(N)T} \mathbf{A}^{(N)})$ ;
7     for  $n = 1$  to  $N$  do
8       if  $n = N$  then
9          $\gamma = 1$ ;
10      end
11       $\mathbf{T}^{(2)} = \mathbf{Y}_{(n)} \{\mathbf{A}^{\odot -n}\}$ ;
12       $\mathbf{T}^{(3)} = \mathbf{T}^{(1)} \oslash (\mathbf{A}^{(n)T} \mathbf{A}^{(n)})$ ;
13      for  $j = 1$  to  $J$  do
14         $\mathbf{a}_j^{(n)} \leftarrow [\gamma_j \mathbf{a}_j^{(n)} + \mathbf{t}_j^{(2)} - \mathbf{A}^{(n)} \mathbf{t}_j^{(3)}]_+$ ;
15        if  $n \neq N$  then
16           $\mathbf{a}_j^{(n)} \leftarrow \mathbf{a}_j^{(n)} / \|\mathbf{a}_j^{(n)}\|_2$ ;      /* normalize to unit length */
17        end
18      end
19       $\mathbf{T}^{(1)} = \mathbf{T}^{(3)} \otimes (\mathbf{A}^{(n)T} \mathbf{A}^{(n)})$ ;
20    end
21  until a stopping criterion is met;
22 end
```

---

## 7 Block Principal Pivoting

Another way to compute a Nonnegative Tensor Factorization is called Block Principal Pivoting, and is an active-set-like method [12] [13].

A tensor  $\underline{\mathbf{Y}}$  of size  $I_1 \times I_2 \times \dots \times I_N$  can be transformed into a matrix  $\mathbf{Y}_{MAT(n)}$  in a process called mode- $n$ -matricization by linearizing all indices except  $n$  in the following way:  $\mathbf{Y}_{MAT(n)}$  is a matrix of size  $I_n \times \prod_{k=1, k \neq n}^N I_k$  and the  $(i_1, \dots, i_N)$ th element of  $\underline{\mathbf{Y}}$  is mapped to the  $(i_n, L)$ th element of  $\mathbf{Y}_{MAT(n)}$  where

$$L = 1 + \sum_{k=1}^N (i_k - 1)L_k$$

and

$$L_k = \prod_{j=1, j \neq n}^{k-1} I_j$$

For a tensor  $\underline{\mathbf{Y}} \in \mathbb{R}^{I_1 \times I_2 \times \dots \times I_N}$ , its approximation  $\hat{\underline{\mathbf{Y}}} = \llbracket \mathbf{A}^{(1)}, \mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)} \rrbracket$  can be written as, for any  $n \in 1, \dots, N$

$$\mathbf{Y}_{MAT(n)} = \mathbf{A}^{(n)} \times (\mathbf{A}^{\odot-n})^T$$

which can be utilized in the ANLS (alternating nonnegativity-constrained least squares) framework. The following subproblems have to be solved:

$$\min_{\mathbf{A}^{(n)} \geq \mathbf{0}} \left\| \mathbf{A}^{\odot-n} \times (\mathbf{A}^{(n)})^T - (\mathbf{Y}_{MAT(n)})^T \right\|_F^2$$

Algorithm 5 shows the Block Principal Pivoting algorithm that can be used to solve this problem. Since the problem was reduced to a matrix problem, any NMF algorithm could be used to compute the factorization, but algorithms that exploit the special properties of the matrices are especially promising. ( $\mathbf{A}^{\odot-n}$  is typically long and thin, while  $(\mathbf{A}^{(n)})^T$  is typically flat and wide.) In the formulation of the Block Principal Pivoting algorithm below,  $\mathbf{x}_{\mathcal{F}_l}$  and  $\mathbf{y}_{\mathcal{G}_l}$  represent the subset of the  $l$ th column of  $\mathbf{X}$  and  $\mathbf{Y}$  indexed by  $\mathcal{F}_l$  and  $\mathcal{G}_l$ , respectively.

---

**Algorithm 5:** Block principal pivoting (from [13])

---

**Input:**  $\mathbf{V} \in \mathbb{R}^{P \times Q}$ ,  $\mathbf{W} \in \mathbb{R}^{P \times L}$

**Output:**  $\mathbf{X} (\in \mathbb{R}^{Q \times L} = \arg(\min_{\mathbf{x} \geq \mathbf{0}} \|\mathbf{V}\mathbf{X} - \mathbf{W}\|_F^2))$

```
1 begin
2   Compute  $\mathbf{V}^T \mathbf{V}$  and  $\mathbf{V}^T \mathbf{W}$ ;
3   Initialize  $\mathcal{F}_l = \emptyset$  and  $\mathcal{G}_l = \{1, \dots, q\}$  for all  $l \in 1, \dots, L$ ;
4    $\mathbf{X} = \mathbf{0}$ ;
5    $\mathbf{Y} = -\mathbf{V}^T \mathbf{W}$ ;
6    $\alpha (\in \mathbb{R}^r) = \mathbf{3}$ ;
7    $\beta (\in \mathbb{R}^r) = (\mathbf{q} + \mathbf{1})$ ;
8   Compute all  $\mathbf{x}_{\mathcal{F}_l}$  using column grouping from  $\mathbf{V}_{\mathcal{F}_l}^T \mathbf{V}_{\mathcal{F}_l} \mathbf{x}_{\mathcal{F}_l} = \mathbf{V}_{\mathcal{F}_l}^T \mathbf{w}_l$ ;
9   Compute all  $\mathbf{y}_{\mathcal{G}_l} = \mathbf{V}_{\mathcal{G}_l}^T (\mathbf{V}_{\mathcal{F}_l} \mathbf{x}_{\mathcal{F}_l} - \mathbf{w}_l)$ ;
   /*  $(\mathbf{x}_{\mathcal{F}_l}, \mathbf{y}_{\mathcal{G}_l})$  is feasible iff  $\mathbf{x}_{\mathcal{F}_l} \geq \mathbf{0}$  and  $\mathbf{y}_{\mathcal{G}_l} \geq \mathbf{0}$  */
10  while any  $(\mathbf{x}_{\mathcal{F}_l}, \mathbf{y}_{\mathcal{G}_l})$  is infeasible do
11    Find the indices of columns in which the solution is infeasible:
       $I = \{j : (\mathbf{x}_{\mathcal{F}_l}, \mathbf{y}_{\mathcal{G}_l}) \text{ is infeasible}\}$ ;
12    for all the  $l \in I$  do
13       $\mathcal{H}_l = \{q \in \mathcal{F}_l : \mathbf{x}_q < 0\} \cup \{q \in \mathcal{G}_l : \mathbf{y}_q < 0\}$ ;
14    end
15    for all the  $l \in I$  with  $|\mathcal{H}_l| < \beta_l$  do
16       $\beta_l = |\mathcal{H}_l|$ ;
17       $\alpha_l = 3$ ;
18       $\hat{\mathcal{H}}_l = \mathcal{H}_l$ ;
19    end
20    for all the  $l \in I$  with  $|\mathcal{H}_l| \geq \beta_l$  and  $\alpha_l \geq 1$  do
21       $\alpha_l = \alpha_l - 1$ ;
22       $\hat{\mathcal{H}}_l = \mathcal{H}_l$ ;
23    end
24    for all the  $l \in I$  with  $|\mathcal{H}_l| \geq \beta_l$  and  $\alpha_l = 0$  do
25       $\hat{\mathcal{H}}_l = \{q : q = \max\{q \in \mathcal{H}_l\}\}$ ;
26    end
27    for all the  $l \in I$  do
28       $\mathcal{F}_l = (\mathcal{F}_l - \hat{\mathcal{H}}_l) \cup (\hat{\mathcal{H}}_l \cap \mathcal{G}_l)$ ;
29       $\mathcal{G}_l = (\mathcal{G}_l - \hat{\mathcal{H}}_l) \cup (\hat{\mathcal{H}}_l \cap \mathcal{F}_l)$ ;
30      Compute  $\mathbf{x}_{\mathcal{F}_l}$  using column grouping from  $\mathbf{V}_{\mathcal{F}_l}^T \mathbf{V}_{\mathcal{F}_l} \mathbf{x}_{\mathcal{F}_l} = \mathbf{V}_{\mathcal{F}_l}^T \mathbf{w}_l$ ;
31       $\mathbf{y}_{\mathcal{G}_l} = \mathbf{V}_{\mathcal{G}_l}^T (\mathbf{V}_{\mathcal{F}_l} \mathbf{x}_{\mathcal{F}_l} - \mathbf{w}_l)$ ;
32    end
33  end
34 end
```

---

## 8 Conclusion

In this report, the Nonnegative Tensor Factorization problem was defined and five algorithms to compute this factorization were presented. For future work, it would be especially interesting to parallelize some of these algorithms to allow the processing of large problems which arise in real-world applications on highly parallel modern computing systems.

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