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TENSOR DECOMPOSITIONS
THEORY AND IMPLEMENTATION

1 CP Decomposition

By convention, let $\mathbf{A} \in \mathbb{R}^{I \times J}$ be a matrix, we shall denote by $\mathbf{a}_j \in \mathbb{R}^I$, where $1 \leq j \leq J$ the j-th column of \mathbf{A} . Let $\mathcal{X} \in \mathbb{R}^{I \times J \times K}$ be an order-3 tensor. The CP decomposition seeks for matrices $\mathbf{A} \in \mathbb{R}^{I \times R}$, $B \in \mathbb{R}^{J \times R}$ and $C \in \mathbb{R}^{K \times R}$ such that

$$\mathcal{X} = \sum_{r=1}^{R} \mathbf{a}_r \circ \mathbf{b}_r \circ \mathbf{c}_r,\tag{1}$$

such that R is the smallest possible number. In such case, R is called the rank of \mathcal{X} . It is usually useful to standardize the column vectors of the matrices to unit norm by introducing scaling factors $\boldsymbol{\lambda} = (\lambda_1, \dots, \lambda_r)^{\top}$, i.e.

$$\mathcal{X} = \sum_{r=1}^{R} \lambda_r \, \mathbf{a}_r \circ \mathbf{b}_r \circ \mathbf{c}_r. \tag{2}$$

The decomposition can be described briefly as $\mathcal{X} = [[\boldsymbol{\lambda}, \mathbf{A}, \mathbf{B}, \mathbf{C}]]$. Let $\Lambda = \operatorname{diag}(\lambda)$, we have the equivalent matricization expressions

$$\begin{cases} \mathbf{X}_{(1)} = \mathbf{A}\Lambda(\mathbf{C}\odot\mathbf{B})^{\top} \\ \mathbf{X}_{(2)} = \mathbf{B}\Lambda(\mathbf{C}\odot\mathbf{A})^{\top} \\ \mathbf{X}_{(3)} = \mathbf{C}\Lambda(\mathbf{B}\odot\mathbf{A})^{\top}. \end{cases}$$
(3)

There is no known finite algorithm for determining the rank of a tensor [1]. Therefore, for each R = 1, 2, ..., we attempt to minimize the difference between \mathcal{X} and a rank-R tensor $\hat{\mathcal{X}} = [[\boldsymbol{\lambda}, \mathbf{A}, \mathbf{B}, \mathbf{C}]]$, i.e.

$$\min_{\hat{\mathcal{X}}} \|\mathcal{X} - \hat{\mathcal{X}}\|^2 \tag{4}$$

until the difference is less than a given tolerance. Using (3), we can write, for example, mode-1 matricization form of the problem as

$$\min_{\hat{\mathcal{Y}}} \| \mathcal{X}_{(1)} - \mathbf{A}\Lambda(\mathbf{C} \odot \mathbf{B}) \|. \tag{5}$$

The solution is

$$\mathbf{A}\Lambda = \mathcal{X}_{(1)}[(\mathbf{C} \odot \mathbf{B})^{\top}]^{\dagger} = \mathcal{X}_{(1)}(\mathbf{C} \odot \mathbf{B})(\mathbf{C}^{\top}\mathbf{C} \star \mathbf{B}^{\top}\mathbf{B})^{\dagger}$$
(6)

We use the last expression because it requires calculating the pseudoinverse of an $R \times R$ matrix rather than a $JK \times R$ matrix. Therefore, we can solve for each factor iteratively until a convergence criterion is met. The general case for N-way tensor is represented in the algorithm below.

Algorithm 1 CP Decomposition using Alternative Least Square

Input:
$$\mathcal{X} \in \mathbb{R}^{I_1,...,I_N}$$
, $\mathbf{A}_{(n)} \in \mathbb{R}^{I_n \times R}$, $\lambda \in \mathbb{R}^R$, $\epsilon = 1e-10$, maxIter = 5000. iter $\leftarrow 0$ while $\|\mathcal{X} - [[\boldsymbol{\lambda}, \mathbf{A}_{(1)}, \ldots, \mathbf{A}_{(n)}]]\| > \epsilon$ and iter $<$ maxIter do for $n = 1, \ldots, N$ do $\mathbf{V} \leftarrow \mathbf{A}_{(1)}^{\top} \mathbf{A}_{(1)} \star \ldots \star \mathbf{A}_{(n-1)}^{\top} \mathbf{A}_{(n-1)} \star \mathbf{A}_{(n+1)}^{\top} \mathbf{A}_{(n+1)} \star \ldots \star \mathbf{A}_{(N)}^{\top} \mathbf{A}_{(N)}$ $\mathbf{A}_{(n)} \leftarrow \mathbf{X}_{(n)} (\mathbf{A}_{(N)} \odot \ldots \odot \mathbf{A}_{(n+1)} \mathbf{A}_{(n-1)} \odot \ldots \odot \mathbf{A}_{(1)}) \mathbf{V}^{\dagger}$ $\lambda \leftarrow (\mathbf{a}_{(n)1}, \ldots, \mathbf{a}_{(n)r})^{\top}$ Normalize columns of $\mathbf{A}_{(N)}$ end for end while Output: $\lambda, \mathbf{A}_{(1)}, \ldots, \mathbf{A}_{(N)}$.

2 Tucker Decompositions: HOSVD and HOOI

Let $R_n = \operatorname{rank}(\mathbf{X}_{(n)})$. Tucker decomposition of ranks S_1, \ldots, S_N , where $S_n \leq R_n$ seeks a tensor $\mathcal{G} \in \mathbb{R}^{S_1 \times \ldots \times S_N}$ and matrices $\mathbf{A}_{(n)} \in \mathbb{R}^{I_n \times S_n}$, where $n = 1, \ldots, N$ to minimize

$$\|\mathcal{X} - \mathcal{G} \times_1 \mathbf{A}_{(1)} \dots \times_N \mathbf{A}_{(N)}\|. \tag{7}$$

The decomposition is written briefly as

$$\mathcal{X} \approx [[\mathcal{G}, \mathbf{A}_{(1)}, \dots, \mathbf{A}_{(n)}]]. \tag{8}$$

If $S_n = R_n$, for all n = 1, ..., N, we have an exact such decomposition, called the higher-order SVD of \mathcal{X} (HOSVD).

Algorithm 2 Higher-order SVD

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Input: \mathcal{X} \in \mathbb{R}^{I_1,...,I_N}, \mathcal{G} \in \mathbb{R}^{R_1,...,R_N}, \mathbf{A}_{(n)} \in \mathbb{R}^{I_n \times R_n}.

for n = 1, ..., N do
\mathbf{A}_{(n)} \leftarrow R_n \text{ leading left singular vectors of } \mathbf{X}_{(n)}
end for
\mathcal{G} = \mathcal{X} \times_1 \mathbf{A}_{(1)}^\top ... \times_N \mathbf{A}_{(N)}^\top
Output: \mathcal{G}, \mathbf{A}_{(1)}, ..., \mathbf{A}_{(N)}.
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In the case there exists $S_n < R_n$ for some n = 1, ..., N, taking S_n left singular vectors of $\mathbf{X}_{(n)}$ as in HOSVD does not lead to an optimal solution. We use the Higher-order Orthogonal Iteration (HOOI) approach with HOSVD as an initialization.

Algorithm 3 Higher-order Orthogonal Iteration

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Input: \mathcal{X} \in \mathbb{R}^{I_1, \dots, I_N}, \mathcal{G} \in \mathbb{R}^{S_1, \dots, S_N}, \mathbf{A}_{(n)} \in \mathbb{R}^{I_n \times S_n}, \epsilon = 1e - 10, maxIter = 5000. Initialize \mathbf{A}_{(n)}, n = 1, \dots, N using HOSVD.

while \|\mathcal{X} - [[\mathcal{G}, \mathbf{A}_{(1)}, \dots, \mathbf{A}_{(n)}]]\| > \epsilon and iter < maxIter do

for n = 1, \dots, N do

\mathcal{Y} \leftarrow \mathcal{X} \times_1 \mathbf{A}_{(1)}^{\top} \dots \times_{n-1} \mathbf{A}_{(n-1)}^{\top} \times_{n+1} \mathbf{A}_{(n+1)}^{\top} \dots \times_N \mathbf{A}_{(N)}^{\top}

\mathbf{A}_{(n)} \leftarrow S_n singular vectors of \mathbf{Y}_{(n)}

end for

end while

\mathcal{G} = \mathcal{X} \times_1 \mathbf{A}_{(1)}^{\top} \dots \times_N \mathbf{A}_{(N)}^{\top}

Output: \mathcal{G}, \mathbf{A}_{(1)}, \dots, \mathbf{A}_{(N)}.
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3 Implementation Details

References

[1] Joseph B Kruskal. "Rank, decomposition, and uniqueness for 3-way and N-way arrays". In: *Multiway data analysis*. 1989, pp. 7–18.