

A Practical Randomized CP Tensor Decomposition

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What is the Paper About?

- CP/PARAFAC Decomposition is a useful tool in data analytics
- The algorithm to compute the decomposition using the Alternating Least Squares Method (ALS) is slow
- Paper attempts to solve this by reducing computation time and memory

How do they do it?

- Randomized methods have been useful in solving Linear Least Squares Problem

Important Equations and Results

- Tensor $\mathcal{X} \in \mathbb{R}^{I_1 \times \cdots \times I_N}$
- mode- n matricization with one-indexing

$$x_{i_1, i_2, \dots, i_N} = \left[\mathbf{X}_{(n)} \right]_{(i_n, j)} \quad (1)$$

$$j = 1 + \sum_{\substack{k=1 \\ k \neq n}}^N (i_k - 1) J_k \quad \text{where } J_k = \prod_{\substack{m=1 \\ m \neq n}}^{k-1} I_m. \quad (2)$$

Important Equations and Results

$\mathbf{A} \in \mathbb{R}^{I \times J}$, $\mathbf{B} \in \mathbb{R}^{I \times J}$

$$(\mathbf{A} \odot \mathbf{B})^T (\mathbf{A} \odot \mathbf{B}) = \mathbf{A}^T \mathbf{A} * \mathbf{B}^T \mathbf{B} \quad (3)$$

$$\mathbf{AB} \otimes \mathbf{CD} = (\mathbf{A} \otimes \mathbf{C})(\mathbf{B} \otimes \mathbf{D}) \quad (4)$$

$$\mathbf{AB} \odot \mathbf{CD} = (\mathbf{A} \otimes \mathbf{C})(\mathbf{B} \odot \mathbf{D}). \quad (5)$$

Important Equations and Results

$$\mathcal{Y} = \mathcal{X} \times_n \mathbf{A} \Leftrightarrow \mathbf{Y}_{(n)} = \mathbf{A} \mathbf{X}_{(n)}. \quad (6)$$

$$\mathcal{Y} = \mathcal{X} \times_1 \mathbf{U}^{(1)} \dots \times_N \mathbf{U}^{(N)} \Leftrightarrow \quad (7)$$

$$\mathbf{Y}^{(n)} = \mathbf{U}^{(n)} \mathbf{X}_{(n)} \left(\mathbf{U}^{(N)} \otimes \dots \otimes \mathbf{U}^{(n+1)} \otimes \mathbf{U}^{(n-1)} \otimes \dots \otimes \mathbf{U}^{(1)} \right)^T. \quad (8)$$

Important Equations and Results

CP decomposition approximates order-N tensor as a sum of R rank-one tensors

$$\mathcal{X} \approx \tilde{\mathcal{X}} = \sum_{r=1}^R a_r^{(1)} \circ a_r^{(2)} \circ \dots \circ a_r^{(N)} \quad (9)$$

$$\mathbf{A}^{(n)} = \begin{bmatrix} \mathbf{a}_1^{(n)} & \mathbf{a}_2^{(n)} & \dots & \mathbf{a}_R^{(n)} \end{bmatrix} \in \mathbb{R}^{I_n \times R} \quad (10)$$

Important Equations and Results

Mode- n matricization in terms of factor matrices

$$\tilde{\mathbf{X}}_{(n)} = \mathbf{A}^{(n)} \mathbf{Z}^{(n)\top} \quad (11)$$

$$\mathbf{Z}^{(n)} = \mathbf{A}^{(N)} \odot \dots \odot \mathbf{A}^{(n+1)} \odot \mathbf{A}^{(n-1)} \odot \dots \odot \mathbf{A}^{(1)} \quad (12)$$

Introducing a normalizing factor λ_r

$$\tilde{\mathcal{X}} = \sum_{r=1}^R \lambda_r \mathbf{a}_r^{(1)} \circ \mathbf{a}_r^{(2)} \circ \dots \circ \mathbf{a}_r^{(N)} \quad (13)$$

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Sketching

$$\min \|Ax - b\| \rightarrow \min \|SAx - Sb\|$$

The leverages of a matrix A are the norms of the rows of the left singular Matrix U

$$l_i = \|U(i, :)^2\|$$

Shows the importance of the row in constructing the column space of the Matrix A

The coherence of a matrix is the maximum leverage scores amongst its rows. It is a measure of how "spread out" the **importance of the rows** is.

$$\mu(A) = \max_{i \in [n]} l_i$$

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The lower the coherence, the better random sketching gets!

$$\mu(A) \geq \frac{d}{n} \quad | A \in \mathbb{R}^{n \times d}$$

- This is done to mix the rows of A such that the coherences can be evenly spread before random sampling.
- Moreover, we use special unitary matrices F (like a DFT Matrix) in order to perform matrix computation fast (using FFT) and a Diagonal Matrix D . (See [DMMS07]) for more details.

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The accuracy of this operation can be provided using FJLT and its variants.

Objective

- Fitting CP method using Alternating Least Squares Method
- Alternate among modes and fix all factor matrices except $\mathbf{A}^{(n)}$
- From Eq. 13, we get the objective as:

$$\arg \min_{\mathbf{A}^{(n)}} \|\mathbf{X}_{(n)} - \mathbf{A}^{(n)} \mathbf{Z}^{(n)T}\|_F. \quad (14)$$

Normal Equation

In CP-ALS, normal equation of Eq. 14 is used

$$\mathbf{X}_{(n)} \mathbf{Z}^{(n)} = \mathbf{A}^{(n)} \left(\mathbf{Z}^{(n)\top} \mathbf{Z}^{(n)} \right) \quad (15)$$

Normal Equation

Using Eq. 3, we get:

$$\mathbf{Z}^{(n)\top} \mathbf{Z}^{(n)} = \mathbf{A}^{(N)\top} \mathbf{A}^{(N)} * \dots * \mathbf{A}^{(n+1)\top} \mathbf{A}^{(n+1)} \\ * \mathbf{A}^{(n-1)\top} \mathbf{A}^{(n-1)} * \dots * \mathbf{A}^{(1)\top} \mathbf{A}^{(1)} \quad (16)$$

Initialization

- $\mathbf{A}^{(n)}$ can be initialized using HOSVD
- It can also be initialized using a random distribution

Algorithm 1: CP-ALS Algorithm

Input: $\mathcal{X} \in \mathbb{R}^{I_1 \times \dots \times I_N}$, \mathbf{R}

Output: λ , $\{\mathbf{A}^{(n)}\}$

```

1 Function CP-ALS( $\mathcal{X}$ ,  $\mathbf{R}$ ):
2   Initialize factor matrices  $\mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)}$ 
3   repeat
4     for  $n = 1, \dots, N$  do
5        $\mathbf{V} \leftarrow \mathbf{A}^{(N)T} \mathbf{A}^{(N)} * \dots * \mathbf{A}^{(n+1)T} \mathbf{A}^{(n+1)} * \mathbf{A}^{(n-1)T} \mathbf{A}^{(n-1)} * \dots * \mathbf{A}^{(1)T} \mathbf{A}^{(1)}$ 
6        $\mathbf{Z}^{(n)} \leftarrow \mathbf{A}^{(N)} \odot \dots \odot \mathbf{A}^{(n+1)} \odot \mathbf{A}^{(n-1)} \odot \dots \odot \mathbf{A}^{(1)}$ 
7        $\mathbf{W} \leftarrow \mathbf{X}^{(n)} \mathbf{Z}^{(n)}$ 
8       Solve  $\mathbf{A}^{(n)} \mathbf{V} = \mathbf{W}$  for  $\mathbf{A}^{(n)}$ 
9       Normalize columns of  $\mathbf{A}^{(n)}$  and update  $\lambda$ 
10    end
11  until termination criteria met
12  return  $\lambda$ , factor matrices  $\{\mathbf{A}^{(n)}\}$ 
  
```

Computation Cost

- Line 5: $O(R^2 \sum_{m \neq n} I_m) + O(R^2 N)$
- Line 6: $O(R \prod_{m \neq n} I_m)$
- Line 7: $O(R \prod_m I_m)$
- Line 8: $O(2R^2 I_n)$
- Outer Iteration: $O(NR \prod_n I_n)$
- Initialization: $(\sum_{n>1} I_n) \prod_m I_m + \sum_{n>1} I_n^3$

Modifying Equations

We will rewrite Eq. 14 as :

$$\arg \min_{\mathbf{A}^{(n)}} \|\mathbf{A}^{(n)} \mathbf{Z}^{(n)T} - \mathbf{X}_{(n)}\|_F. \quad (17)$$

Random Sampling

- Uniformly sample rows from $\mathbf{Z}^{(n)}$ and the corresponding rows from $\mathbf{X}_{(n)}^T$
- Let S denote the number of desired number of samples
- Let \mathcal{S} denote the samples from $\{1, \dots, \prod_{m \neq n} l_m\}$ and $|\mathcal{S}| = S$
- Rows of \mathbf{S} of $\prod_{m \neq n} l_m \times \prod_{m \neq n} l_m$ identity matrix

How is it Better?

- Forming the (full) Khatri-Rao product $\mathbf{Z}^{(n)}$ is expensive
- Compute $\mathbf{SZ}^{(n)}$ without computing $\mathbf{Z}^{(n)}$
- Using matricization mapping, the j^{th} row of $\mathbf{Z}^{(n)}$ of Hadamard Product

$$\mathbf{Z}^{(n)}(j, :) = \mathbf{A}^{(1)}(i_1, :) * \cdots * \mathbf{A}^{(n-1)}(i_{n-1}, :) \\ * \mathbf{A}^{(n+1)}(i_{n+1}, :) * \cdots * \mathbf{A}^{(N)}(i_N, :) \quad (18)$$

Algorithm 2: Sampled Khatri-Rao Product

Input: $\mathbf{S}, \mathbf{A}^{(N)}, \dots, \mathbf{A}^{(n+1)}, \mathbf{A}^{(n-1)}, \dots, \mathbf{A}^{(1)}$
Output: \mathbf{Z}_S

```

1 Function SKR( $\mathbf{S}, \mathbf{A}^{(N)}, \dots, \mathbf{A}^{(n+1)}, \mathbf{A}^{(n-1)}, \dots, \mathbf{A}^{(1)}$ ):
2   Retrieve idxs from  $\mathbf{S}$ 
3    $\mathbf{Z}_S \leftarrow \mathbf{1}$   $\triangleright \mathbf{1} \in \mathbb{R}^{S \times R}$ 
4   for  $m = 1, \dots, n-1, n+1, \dots, N$  do
5      $\mathbf{A}_S^{(m)} \leftarrow \mathbf{A}^{(m)}(\text{idxs}(:, m), :)$   $\triangleright$  MATLAB-style indexing
6      $\mathbf{Z}_S \leftarrow \mathbf{Z}_S * \mathbf{A}_S^{(m)}$ 
7   end
8   return  $\mathbf{Z}_S$ 

```

Algorithm 3: CPRAND Algorithm

Input: $\mathcal{X} \in \mathbb{R}^{I_1 \times \dots \times I_N}$, \mathbf{R} , \mathbf{S}

Output: λ , $\{\mathbf{A}^{(n)}\}$

```

1  Function CPRAND( $\mathcal{X}$ ,  $\mathbf{R}$ ,  $\mathbf{S}$ ):
2      Initialize factor matrices  $\mathbf{A}^{(2)}, \dots, \mathbf{A}^{(N)}$ 
3      repeat
4          for  $n = 1, \dots, N$  do
5              Define sampling operator  $\mathbf{S} \in \mathbb{R}^{S \times \prod_{m \neq n} I_m}$ 
6               $\mathbf{Z}_S \leftarrow \text{SKR}(\mathbf{S}, \mathbf{A}^{(1)}, \dots, \mathbf{A}^{(n-1)}, \mathbf{A}^{(n+1)}, \dots, \mathbf{A}^{(N)})$ 
7               $\mathbf{X}_S^T \leftarrow \mathbf{S} \mathbf{X}_{(n)}^T$ 
8               $\mathbf{A}^{(n)} \leftarrow \arg \min_{\mathbf{A}} \|\mathbf{Z}_S \mathbf{A}^T - \mathbf{X}_S^T\|_F$ 
9              Normalize columns of  $\mathbf{A}^{(n)}$  and update  $\lambda$ 
10         end
11     until termination criteria met
12     return  $\lambda$ , factor matrices  $\{\mathbf{A}^{(n)}\}$ 
  
```

Computation Cost

- Generating S random multi-indices: $O(SN)$
- Line 6: $O(SR(N - 1))$
- Line 8: $O(2SR^2 + 2SRI_n + R^2I_n)$. Assuming $I_n > R$ and $S > R$, we get $O(2SRI_n)$
- Outer Iteration: $O(SR \sum_n I_n)$

Coherence

- Effectiveness of CPRAND depends on the coherence of the coefficient matrix $\mathbf{Z}^{(n)}$

Lemma 3

Given $\mathbf{A} \in \mathbb{R}^{I \times J}$ and $\mathbf{B} \in \mathbb{R}^{K \times L}$, then $\mu(\mathbf{A} \otimes \mathbf{B}) = \mu(\mathbf{A})\mu(\mathbf{B})$

Lemma 4

Given $\mathbf{A} \in \mathbb{R}^{I \times J}$ and $\mathbf{B} \in \mathbb{R}^{K \times L}$, then $\mu(\mathbf{A} \odot \mathbf{B}) \leq \mu(\mathbf{A})\mu(\mathbf{B})$

Failure of CPRAND

- If the individual factor matrices happen to be highly coherent, CPRAND may fail to converge
- This can be avoided by mixing terms before sampling
- Apply FJLT transformation to inner iteration before sampling

$$\arg \min_{\mathbf{A}^{(n)}} \|\mathcal{F} \mathbf{D} \mathbf{Z}^{(n)} \mathbf{A}^{(n)T} - \mathcal{F} \mathbf{D} \mathbf{X}_{(n)}\|_F \quad (19)$$

$$\hat{\mathbf{Z}}^{(n)} = \bigodot_{\substack{m=N \\ m \neq n}}^1 \mathcal{F}_m \mathbf{D}_m \mathbf{A}^{(m)} = \left(\bigotimes_{\substack{m=N \\ m \neq n}}^1 \mathcal{F}_m \mathbf{D}_m \right) \mathbf{Z}^{(n)} = \left(\bigotimes_{\substack{m=N \\ m \neq n}}^1 \mathcal{F}_m \right) \left(\bigotimes_{\substack{m=N \\ m \neq n}}^1 \mathbf{D}_m \right) \mathbf{Z}^{(n)} \quad (20)$$

Equation - Before Sampling

$$\arg \min_{\mathbf{A}^{(n)}} \left\| \left(\bigotimes_{\substack{m=N \\ m \neq n}}^1 \mathcal{F}_m \mathbf{D}_m \right) \mathbf{Z}^{(n)} \mathbf{A}^{(n)T} - \left(\bigotimes_{\substack{m=N \\ m \neq n}}^1 \mathcal{F}_m \mathbf{D}_m \right) \mathbf{X}^{(n)T} \right\|_F \quad (21)$$

Equation - After Sampling

$$\arg \min_{\mathbf{A}^{(n)}} \left\| \mathbf{S} \hat{\mathbf{Z}}^{(n)} \mathbf{A}^{(n)T} - \mathbf{S} (\mathbf{D}_n \mathcal{F}_n \mathbf{X}^{(n)})^T \right\|_F \quad (22)$$

Algorithm 4: CPRAND-MIX Algorithm

Input: $\mathcal{X} \in \mathbb{R}^{I_1 \times \dots \times I_N}$, \mathbf{R} , \mathbf{S} **Output:** λ , $\{\mathbf{A}^{(n)}\}$

```

1  Procedure CPRAND-MIX( $\mathcal{X}$ ,  $\mathbf{R}$ ,  $\mathbf{S}$ ):
2      Initialize factor matrices  $\mathbf{A}^{(m)}$ ,  $m \in \{2, \dots, N\}$ 
3      Define random sign-flip operators  $\mathbf{D}_m$  and unitary matrices  $\mathcal{F}_m$ ,
         $m \in \{1, \dots, N\}$ 
4      Mix factor matrices:  $\hat{\mathbf{A}}^{(m)} \leftarrow \mathcal{F}_m \mathbf{D}_m \mathbf{A}^{(m)}$ ,  $m \in \{2, \dots, N\}$ 
5      Mix tensor:  $\hat{\mathcal{X}} \leftarrow \mathcal{X} \times_1 \mathcal{F}_1 \mathbf{D}_1 \cdots \times_N \mathcal{F}_N \mathbf{D}_N$ 
6      repeat
7          for  $n = 1, \dots, N$  do
8              Define sampling operator  $\mathbf{S} \in \mathbb{R}^{S \times \prod_{m \neq n} I_m}$ 
9               $\hat{\mathbf{Z}}_S \leftarrow \text{SKR}(\mathbf{S}, \hat{\mathbf{A}}^{(N)}, \dots, \hat{\mathbf{A}}^{(n+1)}, \hat{\mathbf{A}}^{(n-1)}, \dots, \hat{\mathbf{A}}^{(1)})$ 
10              $\hat{\mathbf{X}}_S^T \leftarrow \mathbf{D}_n \mathbf{F}_n^* (\mathbf{S} \hat{\mathcal{X}}_{(n)}^T)^T$ 
11              $\mathbf{A}^{(n)} \leftarrow \arg \min_{\mathbf{A}} \|\hat{\mathbf{Z}}_S \mathbf{A}^T - \hat{\mathbf{X}}_S^T\|_F$ , subject to  $\mathbf{A}$  being real-valued
12             Normalize columns of  $\mathbf{A}^{(n)}$  and update  $\lambda$ 
13              $\hat{\mathbf{A}}^{(n)} \leftarrow \mathcal{F}_n \mathbf{D}_n \mathbf{A}^{(n)}$ 
14         end
15     until termination criteria met
16     return  $\lambda$ , factor matrices  $\{\mathbf{A}^{(n)}\}$ 
  
```

Pre-Mixing

Algorithm 5 CPRAND-PREMIX

```

1: function CPRAND-PREMIX( $\mathcal{X}, R, S$ )  $\triangleright \mathcal{X} \in \mathbb{R}^{I_1 \times \dots \times I_N}$ 
2:   Define random sign-flip operators  $\mathbf{D}_m$  and orthogonal matrices  $\mathcal{F}_m$ ,  $m \in \{1, \dots, N\}$ 
3:   Mix:  $\hat{\mathcal{X}} \leftarrow \mathcal{X} \times_1 \mathcal{F}_1 \mathbf{D}_1 \times \dots \times_N \mathcal{F}_N \mathbf{D}_N$ 
4:    $[\lambda, \{\hat{\mathbf{A}}^{(n)}\}] = \text{CPRAND}(\hat{\mathcal{X}}, R, S)$ 
5:   for  $n = 1, \dots, N$  do
6:     Unmix:  $\mathbf{A}^{(n)} = \mathbf{D}_n \mathcal{F}_n^T \hat{\mathbf{A}}^{(n)}$ 
7:   end for
8:   return  $\lambda$ , factor matrices  $\{\mathbf{A}^{(n)}\}$ 
9: end function

```

Figure: CP-Rand Premix

Cost

- Extra Cost compared to Algorithm 3 on *Line 3*:

$$O\left(\sum_{k=1}^N \prod_m [l_m \log l_k]\right) = O\left((\prod_m l_m) \log(\prod_m l_m)\right)$$

Stopping Criteria I

Problem with Residual Norm

- Sampled least squares computations are so inexpensive that checking this stopping condition can take longer than the iteration
- Authors propose a sampling based for computing approximate stopping criterion

Notation

Let $[N]$ denote the set $\{1, \dots, N\}$, Let \hat{P} denote natural number. Let us take

$$\hat{\mathcal{I}} \subset \mathcal{I} \equiv [I_1] \otimes [I_2] \otimes \dots \otimes [I_N]$$

as a random subset of \hat{P} indices of \mathcal{X}

Stopping Criteria II

Notation

Let $\mathcal{E} = \mathcal{X} - \tilde{\mathcal{X}}$, we can see that

$$\|\mathcal{E}\|^2 = \sum_{\mathbf{i} \in \mathcal{I}} \mathbf{e}_{\mathbf{i}}^2 = P\mu$$

$$P = \prod_n I_n$$

$$\mu = \text{mean} \{ \mathbf{e}_{\mathbf{i}}^2 \mid \mathbf{i} \in \mathcal{I} \}$$

Approximation

The next step is to approximate μ with $\hat{\mu}$, which is on the subset $\hat{\mathcal{I}}$

$$\mu = \hat{\mu} \quad \text{where} \quad \hat{\mu} = \text{mean} \left\{ \mathbf{e}_i^2 \mid i \in \hat{\mathcal{I}} \right\}$$

The new residual norm can be estimated as

$$\frac{\|\mathcal{E}\|}{\|\mathcal{X}\|} = \frac{(P\mu)^{1/2}}{\|\mathcal{X}\|} \approx \frac{(P\hat{\mu})^{1/2}}{\|\mathcal{X}\|}$$

Multiplicative Chernoff-Hoeffding Bounds

Let $\mu_{\max} = \max_i (e_i^2)$ be the maximum value. For any $\gamma \in (0, 1)$, we can write the upper and lower tail bound as:

$$\begin{aligned}\Pr\{\hat{\mu} \geq (1 + \gamma)\mu\} &\leq \exp\left(-\frac{2\gamma^2\mu^2\hat{P}}{\mu_{\max}^2}\right) \\ \Pr\{\hat{\mu} \leq (1 - \gamma)\mu\} &\leq \exp\left(-\frac{\gamma^2\mu^2\hat{P}}{\mu_{\max}^2}\right)\end{aligned}\tag{23}$$

Stopping Criteria V

Lemma

For any $\gamma \in (0, 1)$, we can bound the relative difference in the approximated and true error as:

$$\Pr \left\{ \sqrt{1-\gamma} \leq \frac{(P\hat{\mu})^{1/2}}{\|\mathcal{E}\|} \leq \sqrt{1+\gamma} \right\} \leq \exp \left(-2 \frac{\gamma^2 \mu^2 \hat{P}}{\mu_{\max}^2} \right) \quad (24)$$

Computation Cost

- Cost of Computing $\hat{P}\hat{\mu}$: $O(\hat{P}RN)$
- Cost of Computing exact error: $O(R \prod_n l_n)$

Computational Time

Data Generation

- **Synthetic Data**

- 1 Generated factor matrices $\mathbf{A}^{(n)} \in \mathbb{R}^{I_n \times R}$, $n \in \{1, \dots, N\}$
- 2 Computed

$$\mathcal{X}_{\text{true}} = \sum_{r=1}^{R_{\text{true}}} \mathbf{a}_r^{(1)} \circ \dots \circ \mathbf{a}_r^{(N)}$$

- 3 Added noise $\mathcal{N} \in \mathbb{R}^{I_1 \times \dots \times I_N}$ to $\mathcal{X}_{\text{true}}$ to get \mathcal{X}

$$\mathcal{X} = \mathcal{X}_{\text{true}} + \eta \left(\frac{\|\mathcal{X}_{\text{true}}\|}{\|\mathcal{N}\|} \right) \mathcal{N}$$

Data Generation

- **Synthetic Data**

- **Coil Data Set**

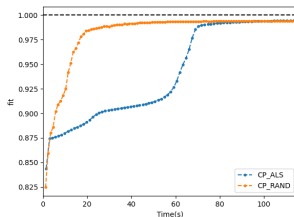
- ① Used the *Coil-100* data set for the experiments.
- ② Contains images of size $128 \times 128 \times 3$ of 100 different objects, with 72 different angles for each object.
- ③ Gives a tensor \mathcal{X} from non Gaussian distribution of the data.
 $\mathcal{X} \in \mathbb{R}^{128 \times 128 \times 3 \times 7200}$

Experiments Conducted

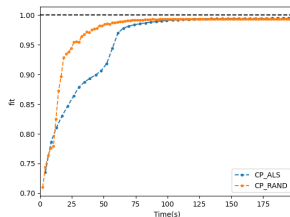
- ① **Fit Time:** Measure the accuracy as a function of time for different models.
- ② **Computational Time:** Measure the average time taken per iteration, ignoring the convergence check.
- ③ **Initialization Time:** Time required to initialize the factor matrices by *Random* and *HOSVD* method.

Experiments

Fit Time



(a) Random $300 \times 300 \times 300$

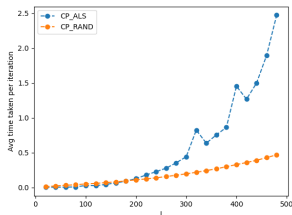


(b) Random $80 \times 80 \times 80 \times 80$

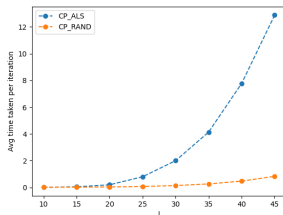
Figure: Runtime comparison for fitting the CP tensor decomposition on random synthetic tensors generated to have rank 5 and 1% noise. We compare a single run of three methods using a target rank of 5

Experiments

Computational Time



(a) order 3: $I \times I \times I$

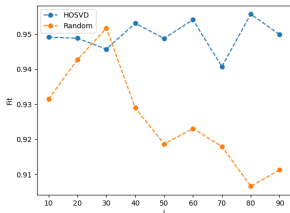


(b) order 5: $I \times I \times I \times I \times I$

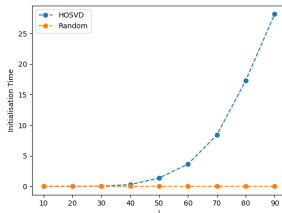
Figure: Mean time per iteration of CP-ALS, CPRAND and CPRAND-MIX for 3rd- and 5th order tensors. The target rank is $R = 5$

Experiments

HOSVD vs Random Initialisation



(a) order 3: $I \times I \times I$



(b) order 3: $I \times I \times I$

Figure: Fig (a) shows the fit value for HOSVD and Random Initialisation for mode 3 tensor, Fig (b) initialisation time for HOSVD and random initialisation, Here data is synthetically generated with actual rank 7, and Target rank 5

Contributions of Paper [BBK17]

- Randomized algorithm prefers incoherent matrices
- Prove that the coherence of the Khatri-Rao product is bounded above by the product of the coherence of its factors
- CPRAND algorithm that uses a randomized least squares solver for the subproblems in CP-ALS
- CPRAND-MIX algorithm that employs efficient mixing to promote incoherence
- Stopping condition that estimates the model fit error



Casey Battaglino, Grey Ballard, and Tamara G. Kolda.
A practical randomized CP tensor decomposition.
CoRR, [abs/1701.06600](https://arxiv.org/abs/1701.06600), 2017.



Petros Drineas, Michael W. Mahoney, S. Muthukrishnan, and
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