

Journées du GDR AMORE, Dec. 2016

Tensor numerical methods for high-dimensional problems

Part 2

Approximation in low-rank formats

We present algorithms for computing low-rank approximations of the solution of **variational problems**

$$\min_{v \in V} \mathcal{J}(v),$$

where V is a tensor space.

- For the approximation of a given tensor u with respect to a certain norm,

$$\mathcal{J}(v) = \|u - v\|.$$

Here, the aim is the compression of u or the extraction of information from u (data analysis).

- For the solution of an equation $Au = b$, the functional $\mathcal{J}(v)$ will measure some distance between u and the approximation v , e.g.

$$\mathcal{J}(v) = \|Av - b\|.$$

The aim is here to obtain an approximation of the solution u with a low computational complexity.

- In **tensor completion**,

$$\mathcal{J}(v) = \sum_{i \in \Omega} |u(i) - v(i)|^2,$$

where $\Omega \subset I$ is a set of known entries of the tensor. The aim is here to **recover (or complete) a tensor from partial information**, by exploiting low-rank structures of the tensor.

- For **inverse problems**, where we want to identify a tensor u from indirect and partial observations, the functional $\mathcal{J}(v)$ measures some distance between observations y and a prediction Av , where A is an observation map:

$$\mathcal{J}(v) = d(y, Av).$$

Exploiting low-rank structures in u allows to reduce the number of parameters to estimate and possibly **makes the problem well-posed**.

- For **least-squares approximation of a function** $u(X)$ from samples $\{u(x^k)\}_{k=1}^n$,

$$\mathcal{J}(v) = \frac{1}{n} \sum_{k=1}^n (u(x^k) - v(x^k))^2$$

- Other problems in statistics and machine learning
 - **Supervised learning** of the relation between a random variable Y and another random variable X from samples $\{(x_k, y_k)\}_{k=1}^n$: minimization of a risk functional

$$\mathcal{J}(v) = \frac{1}{n} \sum_{k=1}^n \ell(y^k, v(x^k))$$

- **Estimation of the density** of a random variable X from samples $\{x_k\}_{k=1}^n$: minimizing the log-likelihood function

$$\mathcal{J}(v) = - \sum_{k=1}^n \log(v(x^k))$$

- ...

- 1 Higher-order singular value decomposition and tensor truncation
- 2 Direct optimization in subsets of low-rank tensors
- 3 Iterative solvers with tensor truncation
- 4 Greedy algorithms

We consider a tensor u in a Hilbert tensor space $V = V^1 \otimes \dots \otimes V^d$ and we assume that u is given as a full tensor or in a certain low-rank format.

We present truncation schemes for finding a low-rank approximation of u with reduced complexity, relying on the standard singular value decomposition of order-two tensors.

We denote by $\|\cdot\|$ the canonical tensor norm on V .

Truncated singular value decomposition for order-two tensors

Let u be an order-two tensor in the Hilbert space $V \otimes W$, where V and W are Hilbert spaces, and let $\|\cdot\|$ denote the canonical norm on $V \otimes W$ (the Frobenius norm for u a matrix).

Let consider a tensor u in $V \otimes W$ with singular value decomposition

$$u = \sum_{k=1}^N \sigma_k v_k \otimes w_k,$$

where the singular values are sorted by decreasing order.

An element of best approximation of u in the set of tensors with rank bounded by r is provided by the [truncated singular value decomposition](#)

$$u_r = \sum_{k=1}^r \sigma_k v_k \otimes w_k,$$

such that

$$\|u - u_r\|^2 = \min_{\text{rank}(v) \leq r} \|u - v\|^2 = \sum_{k=r+1}^N \sigma_k^2.$$

Truncated singular value decomposition for order-two tensors

An approximation u_r with relative precision ϵ , such that

$$\|u - u_r\| \leq \epsilon \|u\|,$$

can be obtained by choosing a rank r such that

$$\sum_{k=r+1}^N \sigma_k^2 \leq \epsilon \sum_{k=1}^N \sigma_k^2.$$

Remark.

The complexity of computing the singular value decomposition of a tensor u is $O(n^3)$ if $\dim(V) = \dim(W) = O(n)$. If u is given in low-rank format $u = \sum_{k=1}^R a_k \otimes b_k$, with a rank $R < n$, the complexity breaks down to $O(R^3 + 2Rn^2)$.

Higher-order tensors as order-two tensors...

For a non-empty subset α in $D = \{1, \dots, d\}$, a tensor $u \in V^1 \otimes \dots \otimes V^d$ can be identified with its matricisation

$$\mathcal{M}_\alpha(u) \in V^\alpha \otimes V^{\alpha^c},$$

an order-two tensor which admits a singular value decomposition

$$\mathcal{M}_\alpha(u) = \sum_{k \geq 1} \sigma_k^\alpha v_k^\alpha \otimes w_k^{\alpha^c}.$$

The set $\sigma^\alpha(u) := \{\sigma_k^\alpha\}_{k \geq 1}$ is called the set of α -singular values of u . The α -rank of u is the number of non-zero α -singular values

$$\text{rank}_\alpha(u) = \|\sigma^\alpha(u)\|_0.$$

Higher-order tensors as order-two tensors...

By sorting the α -singular values by decreasing order, an approximation u_r with α -rank r can be obtained by retaining the r largest singular values, i.e.

$$u_r \text{ such that } \mathcal{M}_\alpha(u_r) = \sum_{k=1}^r \sigma_k^\alpha v_k^\alpha \otimes w_k^{\alpha^c},$$

which satisfies

$$\|u - u_r\|^2 = \min_{\text{rank}_\alpha(v) \leq r} \|u - v\|^2 = \sum_{k>r} (\sigma_k^\alpha)^2.$$

There are 2^{d-1} different binary partitions $\alpha \cup \alpha^c$ of D , to each of which corresponds a singular value decomposition and a way to truncate a higher-order tensor !

Truncation scheme for the approximation in Tucker format

For each $\nu \in \{1, \dots, d\}$, we consider the singular value decomposition of the matricisation $\mathcal{M}_\nu(u)$ of a tensor u

$$\mathcal{M}_\nu(u) = \sum_{k \geq 1} \sigma_k^\nu \mathbf{v}_k^\nu \otimes \mathbf{w}_k^\nu.$$

Let $U_{r_\nu}^\nu = \text{span}\{\mathbf{v}_k^\nu\}_{k=1}^{r_\nu}$ be the subspace of V^ν generated by the r_ν dominant left singular vectors of $\mathcal{M}_\nu(u)$, and by $P_{U_{r_\nu}^\nu}$ the orthogonal projection from V^ν to $U_{r_\nu}^\nu$.

The tensor

$$u_r = (P_{U_{r_1}^1} \otimes \dots \otimes P_{U_{r_d}^d})u$$

is a projection of u onto the **reduced tensor space**

$$U_{r_1}^1 \otimes \dots \otimes U_{r_d}^d$$

and therefore,

$$u_r \in \mathcal{T}_r = \{v \in U^1 \otimes \dots \otimes U^d : U^\nu \subset V^\nu, \dim(U^\nu) = r_\nu, 1 \leq \nu \leq d\}.$$

Higher-order singular value decomposition for Tucker format

The operator

$$\mathcal{P}_{r_\nu}^\nu = \mathcal{M}_\nu^{-1} P_{U_{r_\nu}^\nu} \mathcal{M}_\nu = I \otimes \dots \otimes P_{U_{r_\nu}^\nu} \otimes \dots \otimes I$$

is the orthogonal projection from V onto

$$V^1 \otimes \dots \otimes U_{r_\nu}^\nu \otimes \dots \otimes V^d,$$

which is such that

$$\|u - \mathcal{P}_{r_\nu}^\nu u\| = \min_{\text{rank}_\nu(v) \leq r_\nu} \|u - v\| = \sum_{k \geq r_\nu + 1} (\sigma_k^\nu)^2.$$

The approximation u_r can then be written

$$u_r = \mathcal{P}_{r_1}^1 \dots \mathcal{P}_{r_d}^d u,$$

and satisfies

$$\|u - u_r\|^2 = \sum_{\nu=1}^d \|u - \mathcal{P}_{r_\nu}^\nu u\|^2 = \sum_{\nu=1}^d \min_{\text{rank}_\nu(v) \leq r_\nu} \|u - v\|^2,$$

from which we deduce the quasi-optimality property

$$\|u - u_r\| \leq \sqrt{d} \min_{v \in \mathcal{T}_r} \|u - v\|.$$

Truncation scheme for the approximation in Tucker format

Also, from

$$\|u - u_r\|^2 = \sum_{\nu=1}^d \|u - \mathcal{P}_{r_\nu}^\nu u\|^2 = \sum_{\nu=1}^d \sum_{k_\nu > r_\nu} (\sigma_{k_\nu}^\nu)^2,$$

we deduce that if we select the ranks (r_1, \dots, r_d) such that for each ν

$$\sum_{k_\nu > r_\nu} (\sigma_{k_\nu}^\nu)^2 \leq \frac{\epsilon^2}{d} \sum_{k_\nu \geq 1} (\sigma_{k_\nu}^\nu)^2 = \frac{\epsilon^2}{d} \|u\|^2,$$

then the truncated singular value decomposition $\mathcal{P}_{r_\nu}^\nu u$ has a relative precision ϵ/\sqrt{d} and we finally obtain an approximation u_r with relative precision ϵ ,

$$\|u - u_r\| \leq \epsilon \|u\|.$$

Note that the definition of u_r is independent on the order of the projections $\mathcal{P}_{r_\nu}^\nu$.

Truncation scheme for tree-based tensor formats

For tree-based (hierarchical) low-rank tensor formats

$$\mathcal{T}_r^T = \{v : \text{rank}_\alpha(v) \leq r_\alpha, \alpha \in T\},$$

where T is a subset of a dimension partition tree over $D = \{1, \dots, d\}$, a **higher order singular value decomposition** (also called **hierarchical singular value decomposition**) can also be defined from singular value decompositions of matricisations $\mathcal{M}_\alpha(u)$ of a tensor u .

Truncation scheme for tree-based tensor formats

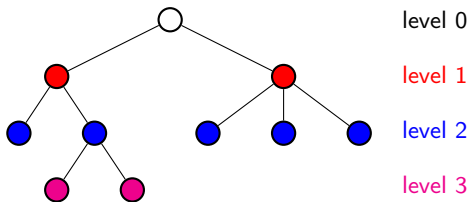
Letting $U_{r_\alpha}^\alpha$ be the subspace generated by the r_α dominant left singular vectors of $\mathcal{M}_\alpha(u)$, and letting $P_{U_{r_\alpha}^\alpha}$ be the orthogonal projector from V^α to $U_{r_\alpha}^\alpha$, we define the orthogonal projection

$$\mathcal{P}_{r_\alpha}^\alpha = \mathcal{M}_\alpha^{-1} P_{U_{r_\alpha}^\alpha} \mathcal{M}_\alpha.$$

Then, an approximation with tree-based rank $r = (r_\alpha)_{\alpha \in T}$ can be defined by

$$u_r = \mathcal{P}_r^{T,(L)} \mathcal{P}_r^{T,(L-1)} \dots \mathcal{P}_r^{T,(1)} u \quad \text{with} \quad \mathcal{P}_r^{T,(\ell)} = \prod_{\substack{\alpha \in T \\ \text{level}(\alpha) = \ell}} \mathcal{P}_{r_\alpha}^\alpha$$

where we apply to u a sequence of projections ordered by increasing level in the tree (from the root to the leaves). Here $L = \max_{\alpha \in T} \text{level}(\alpha)$.



Truncation scheme for tree-based tensor formats

The obtained approximation u_r is such that

$$\|u - u_r\| \leq \sqrt{\#T} \min_{v \in \mathcal{T}_r^T} \|u - v\|,$$

where $\#T \leq 2d - 2$.

Also, if we select the ranks $(r_\alpha)_{\alpha \in T \setminus D}$ such that for all α

$$\sum_{k_\alpha > r_\alpha} (\sigma_{k_\alpha}^\alpha)^2 \leq \frac{\epsilon^2}{\#T} \|u\|^2,$$

we obtain an approximation u_r with relative precision ϵ ,

$$\|u - u_r\| \leq \epsilon \|u\|.$$

Outline

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- 2 Direct optimization in subsets of low-rank tensors
- 3 Iterative solvers with tensor truncation
- 4 Greedy algorithms

Direct optimization in subsets of low-rank tensors

Let \mathcal{M}_r be a subset of tensors in a certain low-rank format \mathcal{M}_r with a **multilinear parametrization** of the form

$$v(i_1, \dots, i_d) = \sum_{k_1=1}^{r_1} \dots \sum_{k_L=1}^{r_L} \prod_{\nu=1}^d p^{(\nu)}(i_\nu, (k_i)_{i \in S_\nu}) \prod_{\nu=d+1}^M p^{(\nu)}((k_i)_{i \in S_\nu})$$

and let

$$\mathcal{M}_r = \{v = \Psi(p^{(1)}, \dots, p^{(M)}) : p^{(\nu)} \in P^{(\nu)}, 1 \leq \nu \leq M\},$$

where Ψ is a multilinear map.

The problem

$$\min_{v \in \mathcal{M}_r} \mathcal{J}(v)$$

can be written as an optimization problem over the parameters

$$\min_{p^{(1)}} \dots \min_{p^{(M)}} \mathcal{J}(\Psi(p^{(1)}, \dots, p^{(M)})).$$

Alternating minimization algorithm

The **alternating minimization algorithm** consists in solving successively minimization problems

$$\min_{\mathbf{p}^{(\nu)} \in \mathcal{P}^{(\nu)}} \mathcal{J}(\Psi(\mathbf{p}^{(1)}, \dots, \mathbf{p}^{(\nu)}, \dots, \mathbf{p}^{(M)})) := \min_{\mathbf{p}^{(\nu)} \in \mathcal{P}^{(\nu)}} \mathcal{J}_\nu(\mathbf{p}^{(\nu)}) \quad (1)$$

over the parameter $\mathbf{p}^{(\nu)}$, letting the other parameters $\mathbf{p}^{(\eta)}$, $\eta \neq \nu$, fixed.

When $\mathcal{P}^{(\nu)}$ is a linear vector space, problem (1) is a **linear approximation problem**.

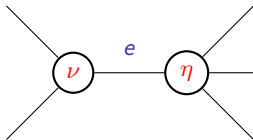
If \mathcal{J} is a **convex** (resp. **differentiable**) functional, then \mathcal{J}_ν is a **convex** (resp. **differentiable**) functional.

Modified alternating minimization algorithm

Modified alternating minimization algorithm¹ is a modification of the alternating minimization algorithm which allows for an **automatic rank adaptation**.

It can be used for optimization in **tree-based tensor formats** or more general **tensor networks**.

At each step of the algorithm, we consider two nodes ν and η connected by an edge e and we update simultaneously the associated parameters $p^{(\nu)}$ and $p^{(\eta)}$.



¹known as **DMRG algorithm** (for Density Matrix Renormalization Group) for tensor networks.

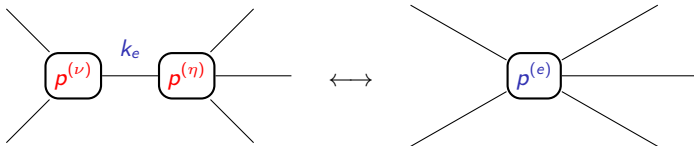
Modified alternating minimization algorithm

In the expression of a tensor $v = \Psi(p^{(1)}, \dots, p^{(M)})$, the two tensors $p^{(\nu)}$ and $p^{(\eta)}$ connected by the edge e appear as

$$\sum_{k_e=1}^{r_e} p^{(\nu)}(k_e, \dots) p^{(\eta)}(k_e, \dots) := p^{(e)}(\dots)$$

where $p^{(e)}$ is a tensor of order

$$\text{order}(p^{(e)}) = \text{order}(p^{(\nu)}) + \text{order}(p^{(\eta)}) - 2.$$



This corresponds to a new tensor networks where the nodes ν and η and edge e are replaced by a single node e , and a new parametrization

$$v = \Psi^e(\dots, p^{(e)}, \dots).$$

Modified alternating minimization algorithm

We first solve an optimization problem

$$\min_{p^{(e)}} \mathcal{J}(\Psi^e(\dots, p^{(e)}, \dots))$$

for obtaining an new value of the tensor $p^{(e)}$.

Then, we compute a low-rank approximation of the tensor $p^{(e)}$

$$p^{(e)}(\dots) \approx \sum_{k_e=1}^{r_e} p^{(\nu)}(k_e, \dots) p^{(\eta)}(k_e, \dots)$$

where the rank r_e in general differs from the initial rank.

In practice, the approximation is obtained using truncated singular value decomposition.

Direct optimization in subsets of low-rank tensors

Other optimization algorithms (e.g. gradient descent, Newton) can be used, possibly exploiting the geometry of low-rank tensor manifolds \mathcal{M}_r .

Under rather standard assumptions, some results have been obtained for the convergence of algorithms: local convergence to a global optimizer, or global convergence to stationary points.

Up to now, there is no available algorithm for obtaining a global optimizer of a general (even convex) functional in a subset of low-rank tensors.

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Iterative solvers with tensor truncation

Another strategy for solving an operator equation

$$Au = b$$

or a more general optimization problem

$$\min_{v \in V} \mathcal{J}(v)$$

is to rely on [classical iterative solvers](#) by interpreting all standard algebraic operations on vector spaces as [algebraic operations in tensor spaces](#).

Iterative solvers with tensor truncation

As a motivating example, consider a simple Richardson algorithm

$$u^n = u^{n-1} - \omega(Au^{n-1} - b).$$

For A and b given in low-rank formats, computing u^n involves **standard algebraic operations**.

However, **the representation rank of the iterates dramatically increases** since

$$\text{rank}(u^n) \approx \text{rank}(A) \text{rank}(u^{n-1}) + \text{rank}(u^{n-1}) + \text{rank}(b).$$

This requires additional **truncation steps for reducing the ranks** of the iterates, such as

$$u^n = T(u^{n-1} - \omega(Au^{n-1} - b)),$$

where $T(v)$ provides a low-rank approximation of v .

We now analyze the behavior of these algorithms depending on the **properties of the truncation operator T** .

Fixed point iterations algorithm

Let us consider a problem which can be written as a fixed point problem

$$F(u) = u,$$

where $F : V \rightarrow V$ is a contractive map, such that for all $u, v \in V$,

$$\|F(u) - F(v)\| \leq \rho \|u - v\|,$$

with $0 \leq \rho < 1$.

Then, consider the fixed point iterations algorithm

$$u^{n+1} = F(u^n)$$

which provides a sequence $(u^n)_{n \geq 1}$ which converges to u , such that

$$\|u - u^n\| \leq \rho^n \|u - u^0\|.$$

Example 1

For a problem $Au = b$, consider $F(u) = u - \omega(Au - b)$, with ω such that $\|I - \omega A\| < 1$. Fixed point iterations $u^{n+1} = u^n - \omega(Au^n - b)$ correspond to Richardson iterations.

Perturbed fixed point iterations algorithm

Now consider the perturbed fixed point iterations

$$v^{n+1} = F(u^n), \quad u^{n+1} = T(v^{n+1})$$

where T is a mapping which for a tensor v provides an **approximation (called truncation)** $T(v)$ in a certain low-rank format \mathcal{M}_r .

Truncations with controlled relative precision

Suppose that the mapping T provides an **approximation with relative precision** ϵ , i.e.

$$\|T(v) - v\| \leq \epsilon \|v\|.$$

This is made possible by using an adaptation of the ranks.

Then the sequence $(u^n)_{n \geq 1}$ is such that

$$\|u - u^n\| \leq \gamma^n \|u - u^0\| + \frac{\epsilon}{1 - \gamma} \|u\|,$$

with $\gamma = \rho(1 + \epsilon)$. Therefore, if $\gamma < 1$

$$\limsup_{n \rightarrow \infty} \|u - u^n\| \leq \frac{\epsilon}{1 - \gamma} \|u\|$$

which means that the sequence tends to **enter a neighborhood of u with radius** $\frac{\epsilon}{1 - \gamma} \|u\|$.

The drawback of this algorithm is that the **ranks of the iterates are not controlled** and may become very high during the iterations.

Truncations in fixed subsets

Now consider that the mapping T provides an approximation in a fixed subset of tensors \mathcal{M}_r with rank bounded by r .

Let us assume that for all v , $T(v)$ provides a quasi-optimal approximation of v such that

$$\|T(v) - v\| \leq C \min_{w \in \mathcal{M}_r} \|v - w\|. \quad (2)$$

A practical realization of a mapping T verifying (2) is provided by [truncated higher-order singular value decompositions](#), where

$$C = O(\sqrt{d}).$$

Truncations in fixed subsets

Let u_r be an element of best approximation of u , with

$$\|u - u_r\| = \min_{v \in \mathcal{M}_r} \|u - v\|.$$

The sequence $(u^n)_{n \geq 1}$ is such that

$$\|u - u^n\| \leq \gamma^n \|u - u^0\| + \frac{C}{1 - \gamma} \|u - u_r\|,$$

with $\gamma = \rho(1 + C)$. If $\gamma < 1$ (which may be quite restrictive on ρ), we obtain

$$\limsup_{n \rightarrow \infty} \|u - u^n\| \leq \frac{C}{1 - \gamma} \min_{v \in \mathcal{M}_r} \|u - v\|,$$

which means that the sequence tends to **enter a neighborhood of u** with radius $\frac{C}{1 - \gamma} \sigma_r$, where σ_r is the best approximation error of u by elements of \mathcal{M}_r .

An advantage of this approach is that the **ranks of the iterates are controlled**. A drawback is that the condition $\gamma < 1$ **imposes to rely on an iterative solver with small contractivity constant** $\rho < (1 + C)^{-1}$, which may be quite restrictive (requires good **preconditioners**).

Truncations with non-expansive maps

Now we assume that the mapping T providing an approximation in low-rank format is non-expansive, i.e.

$$\|T(v) - T(w)\| \leq \|v - w\| \quad (3)$$

The sequence u^n is defined by

$$u^{n+1} = G(u^n),$$

where $G = T \circ F$ is a contractive mapping with the same contractivity constant ρ as F . Therefore, the sequence u^n converges to the unique fixed point u^* of G such that

$$G(u^*) = u^*,$$

with

$$\|u^* - u^n\| \leq \rho^n \|u^* - u^0\|.$$

The obtained approximation u^* is such that

$$(1 + \rho)^{-1} \|u - T(u)\| \leq \|u - u^*\| \leq (1 - \rho)^{-1} \|u - T(u)\|.$$

A practical realization of a mapping T verifying (2) is provided by the **soft singular values thresholding operator**. The **ranks of the iterates are not controlled**. However, it is observed in practice that the **ranks of iterates are usually lower** than with truncations with controlled relative precision.

Convex relaxation

Consider an order two tensor u in a Hilbert tensor space $V \otimes W$ equipped with the canonical norm.

The constrained optimization problem

$$\min_{\text{rank}(v) \leq r} \mathcal{J}(v)$$

is equivalent to

$$\min_v \mathcal{J}(v) + \tau \text{rank}(v)$$

for some value of τ .

A convex optimization problem is obtained by replacing $\text{rank}(v) = \|\sigma(v)\|_0$ by the function $\|\sigma(v)\|_1 = \|v\|_*$ (the trace norm of v)

$$\min_v \mathcal{J}(v) + \tau \|v\|_*$$

Soft thresholding of singular values

Let $u = \sum_{k \geq 1} \sigma_k v_k \otimes w_k$ be the singular value decomposition of u .

The solution of

$$\min_v \frac{1}{2} \|u - v\|^2 + \tau \|v\|_*$$

is given by

$$T_\tau(u) = \sum_{k \geq 1} (\sigma_k - \tau)_+ v_k \otimes w_k$$

which corresponds to a **soft-thresholding of singular values**.

In convex analysis, $T_\tau(u)$ is known as the **proximal operator** of the convex function $v \mapsto \tau \|v\|_*$.

The function $v \mapsto T_\tau(v)$ is non expansive:

$$\|T_\tau(u) - T_\tau(v)\| \leq \|u - v\|.$$

Proximal algorithms

Consider the problem

$$\min_v \mathcal{J}(v) + \tau \|v\|_*$$

A proximal algorithm constructs a sequence $(u^n)_{n \geq 1}$ as follows.

At iteration n , we linearize the function \mathcal{J} around u^n and define u^{n+1} as the solution of

$$\min_v \mathcal{J}(u^n) + (\nabla \mathcal{J}(u^n), v - u^n) + \frac{\beta}{2} \|u - u^n\|^2 + \tau \|v\|_*$$

where β is a parameter.

This is equivalent to solving

$$\min_v \frac{1}{2} \|v - (u^n - \beta^{-1} \nabla \mathcal{J}(u^n))\|^2 + \frac{\tau}{\beta} \|v\|_*$$

whose solution is provided by

$$u^{n+1} = T_{\tau/\beta}(u^n - \beta^{-1} \nabla \mathcal{J}(u^n))$$

where $T_{\tau/\beta}$ is the proximal operator of $v \mapsto \frac{\tau}{\beta} \|v\|_*$.

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 - Greedy algorithms for Tucker format
 - Partially greedy algorithms for Tucker format

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Greedy algorithms for canonical format

A tensor $v \in \mathcal{R}_r$ with **canonical rank** r can be written as a sum of r rank-one tensors

$$v = \sum_{k=1}^r c_k w_k, \quad w_k \in \mathcal{R}_1.$$

Therefore, v can be interpreted as a **n -sparse element with respect to dictionary of rank-one tensors \mathcal{R}_1** .

Greedy algorithms for canonical format

Standard greedy algorithms can be used to construct a **sequence of approximations u^n with increasing canonical rank**

$$u^n = \sum_{k=1}^n c_k^n w_k, \quad c_k^n \in \mathbb{R},$$

where

$$w_n = w_n^{(1)} \otimes \dots \otimes w_n^{(d)} \in \mathcal{R}_1$$

is such that

$$w_n \in \arg \min_{w \in \mathcal{R}_1} \mathcal{J}(u^{n-1} + w), \quad (4)$$

and where the coefficients c_k^n can be either taken as $c_k^n = 1$ (for a pure greedy algorithm), or as the solution of

$$\min_{c_1, \dots, c_n} \mathcal{J}\left(\sum_{k=1}^n c_k w_k\right). \quad (5)$$

Each step requires to solve an **optimization problem in \mathcal{R}_1** , for which we can rely on an alternating minimization algorithm or other optimization algorithms.

Greedy algorithms with dictionary of low-rank tensors

These algorithms are essentially used for the approximation in canonical format but \mathcal{R}_1 could be replaced by another subset of low-rank tensors \mathcal{M} containing \mathcal{R}_1 .

Convergence is guaranteed under quite general assumptions on \mathcal{J} (strongly convex, differentiable with Lipschitz differential) and the set \mathcal{M} (\mathcal{M} closed, $\text{span } \mathcal{M} = V$).

Greedy algorithms with a dictionary \mathcal{R}_1 of rank-one tensors often present a slow convergence compared to the ideal performance of n -term approximations

$$\inf_{v \in \mathcal{R}_n} \mathcal{J}(v).$$

Also, these algorithms do not really exploit the structure of tensors.

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Approximation in Tucker format: a subspace point of view

The set \mathcal{T}_r of tensors with Tucker rank bounded by $r = (r_1, \dots, r_d)$ is defined by

$$\mathcal{T}_r = \left\{ v = \sum_{1 \leq k_1 \leq r_1} \dots \sum_{1 \leq k_d \leq r_d} C_{k_1, \dots, k_d} v_{k_1}^{(1)} \otimes \dots \otimes v_{k_d}^{(d)} : C \in \mathbb{R}^{r_1 \times \dots \times r_d}, v_{k_\nu}^{(\nu)} \in V_\nu \right\}.$$

It can be equivalently parametrized by subspaces

$$\mathcal{T}_r = \{v : v \in U_1 \otimes \dots \otimes U_d \text{ with } U_\nu \subset V_\nu, \dim(U_\nu) = r_\nu\}.$$

Then, an optimization problem on \mathcal{T}_r can be interpreted as a problem of finding **optimal low-dimensional spaces**:

$$\min_{v \in \mathcal{T}_r} \mathcal{J}(v) = \min_{\dim(U_1)=r_1} \dots \min_{\dim(U_d)=r_d} \min_{v \in U_1 \otimes \dots \otimes U_d} \mathcal{J}(v).$$

This is a **multilinear version of projection-based model order-reduction methods**, where an approximation is searched in a tensor product $U_1^{r_1} \otimes \dots \otimes U_d^{r_d}$ of optimal subspaces $U_\nu^{r_\nu}$ of dimension r_ν .

Greedy algorithms for approximation in Tucker format

Greedy algorithms with a [subspace point of view](#), which are similar to greedy algorithms for reduced basis methods, can be introduced for the construction of approximations u^n in an [increasing sequence of tensor subspaces](#)

$$U_1^n \otimes \dots \otimes U_d^n, \quad n \geq 1,$$

with

$$U_\nu^1 \subset \dots \subset U_\nu^n \subset \dots, \quad 1 \leq \nu \leq d.$$

Greedy algorithms for approximation in Tucker format

At step n of these algorithms, we have an approximation u^{n-1} and associated subspaces U_ν^{n-1} of dimension r_ν^{n-1} , $1 \leq \nu \leq d$.

Assume that we have selected a set of dimensions $D_n \subset \{1, \dots, d\}$ to be enriched ($D_n = \{1, \dots, d\}$ for an isotropic enrichment).

For $\nu \notin D_n$, we let $U_\nu^n = U_\nu^{n-1}$, and for $\nu \in D_n$ we construct new spaces U_ν^n with dimension $r_\nu^n = r_\nu^{n-1} + 1$ and such that $U_\nu^n \supset U_\nu^{n-1}$.

An **optimal greedy algorithm** would consist in solving

$$\mathcal{J}(u^n) = \min_{\substack{\dim(U_\nu^n) = r_\nu^n \\ U_\nu^n \supset U_\nu^{n-1} \\ \nu \in D_n}} \min_{v \in U_1^n \otimes \dots \otimes U_d^n} \mathcal{J}(v)$$

Greedy algorithms for approximation in Tucker format

A **practical greedy algorithm** consists in computing an optimal rank-one correction of u^{n-1}

$$\mathcal{J}(u^{n-1} + w_n^{(1)} \otimes \dots \otimes w_n^{(d)}) = \min_{w \in \mathcal{R}_1} \mathcal{J}(u^{n-1} + w),$$

in enriching the spaces according to

$$U_\nu^n = U_\nu^{n-1} + \text{span}(w_n^{(\nu)}), \quad \nu \in D_n,$$

and finally in computing the best approximation u^n in the tensor space $U_1^n \otimes \dots \otimes U_d^n$ by solving

$$\mathcal{J}(u^n) = \min_{v \in U_1^n \otimes \dots \otimes U_d^n} \mathcal{J}(v)$$

or

$$\min_{C \in \mathbb{R}^{r_1^n \times \dots \times r_d^n}} \mathcal{J}\left(\sum_{1 \leq k_1 \leq r_1^n} \dots \sum_{1 \leq k_d \leq r_d^n} C_{k_1 \dots k_d} v_{k_1}^{(1)} \otimes \dots \otimes v_{k_d}^{(d)}\right) \quad (6)$$

where $\{v_i^{(\nu)}\}_{i=1}^{r_\nu^n}$ is a basis of U_ν^n .

For high-dimensional problems, the practical solution of (6) requires a structured approximation of the tensor C , e.g. using sparse or low-rank formats. Note that if we add the constraint of having a super-diagonal tensor C , we recover a standard greedy algorithm for approximation in canonical format.

- 1 Higher-order singular value decomposition and tensor truncation
- 2 Direct optimization in subsets of low-rank tensors
- 3 Iterative solvers with tensor truncation
- 4 Greedy algorithms
 - Greedy algorithms for canonical format
 - Greedy algorithms for Tucker format
 - Partially greedy algorithms for Tucker format

Partially greedy algorithms for Tucker format

For order-two tensors in $V_1 \otimes V_2$, greedy algorithms for Tucker format construct a sequence of spaces

$$U^n = U_1^n \otimes U_2^n,$$

with a greedy enrichment of both left and right spaces, and a corresponding sequence of rank- n approximations u^n with

$$\mathcal{J}(u^n) = \min_{v \in U_1^n \otimes U_2^n} \mathcal{J}(v) = \min_{C \in \mathbb{R}^{n \times n}} \mathcal{J}\left(\sum_{i,j=1}^n v_i^{(1)} \otimes v_j^{(2)} C_{i,j}\right)$$

A **partially greedy strategy** consists in constructing a sequence of spaces

$$U^n = U_1^n \otimes V_2,$$

where only the left spaces are constructed in a greedy fashion.

Partially greedy algorithms for Tucker format

At step n , a suboptimal algorithm consists in computing a rank-one correction of u^{n-1}

$$\mathcal{J}(u^{n-1} + w_n^{(1)} \otimes w_n^{(2)}) = \min_{w^{(1)}, w^{(2)}} \mathcal{J}(u^{n-1} + w^{(1)} \otimes w^{(2)}),$$

in enriching the left subspace according to

$$U_1^n = U_1^{n-1} + \text{span}(w_n^{(1)}),$$

and then in computing an approximation u^n in $U_1^n \otimes V_2$ by solving

$$\mathcal{J}(u^n) = \min_{v \in U_1^n \otimes V_2} \mathcal{J}(v) = \min_{v_1^{(2)}, \dots, v_n^{(2)}} \mathcal{J}\left(\sum_{i=1}^n v_i^{(1)} \otimes v_i^{(2)}\right)$$

where $\{v_i^{(1)}\}_{i=1}^n$ is a basis of U_1^n .

Other topics

- Approximation power of low-rank formats
- Interpolation methods for low-rank approximation
- Geometry of low-rank formats and its consequences in model order reduction of dynamical systems and optimization.
- Selection of a tensor format
- Exploiting sparsity in tensor representations
- Higher-order tensor methods for low-dimensional problems : quantization
- ...