

Structured Low-Rank Approximation and (Some of) Its Applications

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System realisation

The sequence

$$h := (h(0), h(1), \dots), \quad h(t) \in \mathbb{R}^{p \times m}$$

is realisable by a finite dim. LTI system, if and only if

$$\mathcal{H}(h) := \begin{bmatrix} h(1) & h(2) & h(3) & \dots \\ h(2) & h(3) & \ddots & \\ h(3) & \ddots & & \\ \vdots & & & \end{bmatrix}$$

has finite rank. Moreover,

$$\begin{aligned} \text{rank}(\mathcal{H}(h)) &= \text{state dim. of a minimal realisation of } h \\ &= \text{complexity of an exact LTI model for } h. \end{aligned}$$

Examples

Let's start with the following well known examples:

- System realisation
- Discrete deconvolution
- Greatest common divisor of two polynomials

Approximate realisation = Model reduction

However, rank deficiency is a nongeneric property (in $\mathbb{Z}_+ \rightarrow \mathbb{R}^{p \times m}$).

Rank is computed numerically most reliably by the SVD.

From a system theoretic point of view

the SVD does model reduction (Kung's algorithm).

The truncated SVD gives (2-norm) optimal **unstructured** approx.

Instead, we are aiming at a

structured rank- n approximation of $\mathcal{H}(h)$:

Find \hat{h} , such that $\|h - \hat{h}\|$ is minimized and $\text{rank}(\mathcal{H}(\hat{h})) = n$.

Approximate realisation (model reduction)



Hankel structured low-rank approximation

The approximate realisation (model reduction) problem is

Given $h := (h(0), h(1), \dots)$ and $n \in \mathbb{N}$, find

$$\min_{\hat{h}} \|h - \hat{h}\| \quad \text{subject to} \quad \text{rank}(\mathcal{H}(\hat{h})) \leq n$$

a Hankel structured low-rank approximation (SLRA) problem.

Unfortunately, this problem is NP-complete.

Exact and approximate deconvolution

Exact deconv. problem: Given u and y , find h , such that $y = h \star u$.

Solution exists if and only if the system of equations

$$\text{row}(y) = \text{row}(h) \mathcal{T}_{n+1}(u)$$

is solvable for h . However with $T > (n+1)m$, generically solution does not exist \rightsquigarrow **approximate deconvolution problem:**

Given u, y , and $n \in \mathbb{N}$, find

$$\min_{\hat{u}, \hat{y}, \hat{h}} \|\text{col}(u, y) - \text{col}(\hat{u}, \hat{y})\| \quad \text{subject to} \quad \text{row}(\hat{y}) = \text{row}(\hat{h}) \mathcal{T}_{n+1}(\hat{u})$$

Deconvolution

Consider the finite sequences

$$h := (h(0), h(1), \dots, h(n)), \quad \text{where } h \in \mathbb{R}^{p \times m}$$

$$u := (u(-n), \dots, u(0), u(1), \dots, u(T)) \quad \text{and} \quad y := (y(1), \dots, y(T)).$$

Define $\text{row}(y) := [y(1) \quad \dots \quad y(T)]$ and the Toeplitz matrix

$$\mathcal{T}_{n+1}(u) := \begin{bmatrix} u(1) & u(2) & u(3) & \dots & u(T) \\ u(0) & u(1) & u(2) & \dots & u(T-1) \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ u(-n) & u(1-n) & u(2-n) & \dots & u(T-n) \end{bmatrix}$$

With this notation,

$$y = h \star u \quad (\text{convolution}) \quad \Longleftrightarrow \quad \text{row}(y) = \text{row}(h) \mathcal{T}_{n+1}(u) \quad (\text{linear algebra})$$

Deconvolution = FIR system identification

We can interpret

$$y = h \star u$$

as the response of an FIR system with impulse response h to

- initial conditions $(u(-n), \dots, u(0))$, and
- input $(u(1), \dots, u(T))$.

Then the deconvolution problem has the meaning of an **FIR system identification problem:**

Given initial condition, input, and output, find an FIR model.

- exact deconvolution \implies exact FIR fitting model
- approx. deconvolution \implies approx. FIR fitting model

The parameter n bounds the FIR model complexity.

Approximate deconvolution \rightsquigarrow SLRA

Assuming that $\mathcal{T}_{n+1}(\hat{u})$ is full rank (persistency of excitation),

$$\text{row}(\hat{y}) = \text{row}(\hat{h})\mathcal{T}_{n+1}(\hat{u}) \iff \text{rank} \left(\begin{bmatrix} \mathcal{T}_{n+1}(\hat{u}) \\ \text{row}(\hat{y}) \end{bmatrix} \right) = (n+1)m$$

Then the approximate deconvolution problem can be written as

Given u, y , and $n \in \mathbb{N}$, find

$$\min_{\hat{u}, \hat{y}} \|\text{col}(u, y) - \text{col}(\hat{u}, \hat{y})\| \quad \text{subject to}$$

$$\text{rank} \left(\begin{bmatrix} \mathcal{T}_{n+1}(\hat{u}) \\ \text{row}(\hat{y}) \end{bmatrix} \right) \leq (n+1)m$$

a SLRA problem with structure composed of two blocks:
Toeplitz block above an unstructured block.

Approximate GCD \iff Sylvester SLRA

Given $a(z), b(z)$, and $n \in \mathbb{N}$, find

$$\min_{\hat{a}, \hat{b}} \|\text{col}(a, b) - \text{col}(\hat{a}, \hat{b})\| \quad \text{subject to}$$

$$\text{rank}(S(a, b)) \leq n_a + n_b - n$$

Greatest common divisor (GCD)

Consider the polynomials

$$a(z) := a_0 + a_1 z + \dots + a_{n_a} z^{n_a}, \quad b(z) := b_0 + b_1 z + \dots + b_{n_b} z^{n_b}$$

and define the Sylvester matrix

$$S(a, b) := \begin{bmatrix} a_0 & & & b_0 & & \\ \vdots & \ddots & & \vdots & \ddots & \\ a_{n_a} & & a_0 & b_{n_b} & & b_0 \\ & \ddots & \vdots & & \ddots & \vdots \\ & & a_{n_a} & & & b_{n_b} \end{bmatrix} \in \mathbb{R}^{(n_a+n_b) \times (n_a+n_b)}$$

The GCD of $a(z)$ and $b(z)$, has degree n , if and only if

$$\text{rank}(S(a, b)) = n_a + n_b - n.$$

Data matrix being low-rank

an exact property
holds on the data



a matrix constructed
from data is low-rank

- h is realisable by an LTI system of order n $\iff \text{rank}(\mathcal{H}(h)) \leq n$
- (u, y) is fitted by an n taps FIR system $\iff \text{rank} \left(\begin{bmatrix} \mathcal{T}_{n+1}(u) \\ \text{row}(y) \end{bmatrix} \right) \leq (n+1)m$
- $a(z), b(z)$ have GCD of deg. $\geq n$ $\iff \text{rank}(S(a, b)) \leq n_a + n_b - n$

Rank of the data matrix

complexity of an exact
model fitting the data

\leftrightarrow

rank of the
data matrix

- order of the realization $= \text{rank}(\mathcal{H}(h))$
- number of taps
of an FIR system $= \text{rank} \left(\begin{bmatrix} \mathcal{I}_{n+1}(u) \\ \text{row}(y) \end{bmatrix} \right) / m - 1$
- degree of the GCD $= \text{rank deficiency of } S(a, b)$

Structured low-rank approximation

Given

- a vector $p \in \mathbb{R}^{n_p}$,
- a mapping $\mathcal{S} : \mathbb{R}^{n_p} \rightarrow \mathbb{R}^{m \times n}$ (structure specification)
- a vector norm $\|\cdot\|$, and
- an integer r , $0 < r < \min(m, n)$,

find

$$\hat{p}^* := \arg \min_{\hat{p}} \|p - \hat{p}\| \quad \text{subject to} \quad \text{rank}(\mathcal{S}(\hat{p})) \leq r. \quad (*)$$

Interpretation:

$\hat{D}^* := \mathcal{S}(\hat{p}^*)$ is optimal rank- r (or less) approx. of $D := \mathcal{S}(p)$,
within the class of matrices with the same structure as D .

Main issue: Low-rank approximation

With a bounding on the model complexity,

generically in the data space, exact property does not hold

\Rightarrow an approximation is needed.

Approximation paradigm:

modify the data as little as possible, so that the exact property holds for the modified data.

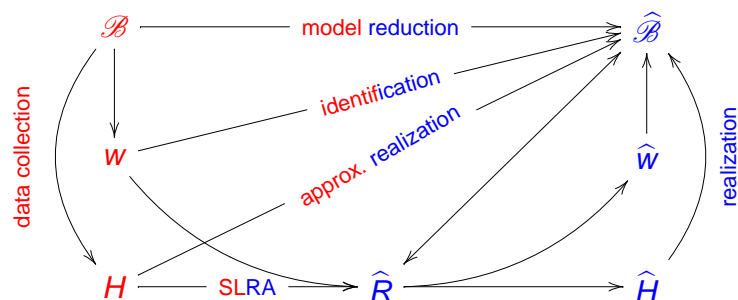
This paradigm leads to structured low-rank approximation.

Applications

- System theory
 1. Approximate realization
 2. Model reduction
 3. Errors-in-variables system identification
 4. Output error system identification
- Signal processing
 5. Output only (autonomous) system identification
 6. Finite impulse response (FIR) system identification
 7. Harmonic retrieval
 8. Image deblurring
- Computer algebra
 9. Approximate greatest common divisor (GCD)

System theory applications

\mathcal{B} “true” (high order) model
 $\hat{\mathcal{B}}$ approximate (low order) model
 w observed response
 H observed impulse resp.
 \hat{w} response of $\hat{\mathcal{B}}$
 \hat{H} impulse resp. of $\hat{\mathcal{B}}$



Statistical vs. deterministic formulation

The EIV model gives a **quality certificate** to the method.

The method works “well” (**consistency**) and is optimal (**efficiency**) under certain specified conditions.

However, the assumption that the data is generated by a true model with additive noise is sometimes not realistic.

Model-data mismatch is often due to a restrictive (LTI) model class being used and not (only) due to measurement noise.

⇒ **The approximation aspect is often more important than the stochastic estimation one.**

Errors-in-variables (EIV) identification

$\mathcal{L}_{m,1}$ LTI model class of bounded complexity ($\#inputs \leq m$, $lag \leq 1$)

Given $w_d \in (\mathbb{R}^w)^T$ and complexity specification $(m, 1)$, find

$$\hat{\mathcal{B}}^* := \arg \min_{\hat{\mathcal{B}}, \hat{w}} \|w_d - \hat{w}\|_{\ell_2} \quad \text{subject to} \quad \hat{w} \in \hat{\mathcal{B}} \in \mathcal{L}_{m,1}.$$

SLRA (*) with $\mathcal{S}(p) = \mathcal{H}_{1+1}(w_d)$, and $r = p$.

EIV model: $w_d = \bar{w} + \tilde{w}$, $\bar{w} \in \bar{\mathcal{B}} \in \mathcal{L}_{m,1}$, $\tilde{w} \sim \text{Normal}(0, \sigma^2 I)$

\bar{w} — true data, $\bar{\mathcal{B}}$ — true model, \tilde{w} — measurement noise

$\hat{\mathcal{B}}^*$ is a maximum likelihood estimate of $\bar{\mathcal{B}}$

consistent and assympt. normal \implies **confidence regions**

Outline

Introduction

Applications

Algorithms

Related problems

Unstructured low-rank approximation

$$\hat{D}^* := \arg \min_{\hat{D}} \|D - \hat{D}\|_F \quad \text{subject to} \quad \text{rank}(\hat{D}) \leq r$$

Theorem (closed form solution)

Let $D = U\Sigma V^\top$ be the SVD of D and define

$$U = \begin{bmatrix} U_1 & U_2 \end{bmatrix} \begin{matrix} r & n-r \\ m & \end{matrix}, \quad \Sigma = \begin{bmatrix} \Sigma_1 & 0 \\ 0 & \Sigma_2 \end{bmatrix} \begin{matrix} r & n-r \\ n-r & \end{matrix} \quad \text{and} \quad V = \begin{bmatrix} V_1 & V_2 \end{bmatrix} \begin{matrix} r & n-r \\ m & \end{matrix}.$$

An optimal low-rank approximation solution is

$$\hat{D}^* = U_1 \Sigma_1 V_1^\top, \quad (\hat{\mathcal{B}}^* = \ker(U_2^\top) = \text{colspan}(U_1)).$$

It is unique if and only if $\sigma_r \neq \sigma_{r+1}$.

Variable projection vs. alternating projections

Two ways to approach the double minimization:

- **Variable projections (VARPRO):**
solve the inner minimization analytically

$$\min_{R, RR^\top = I_{m-r}} \text{vec}^\top(R\mathcal{J}(\hat{p})) \left(G(R)G^\top(R) \right)^{-1} \text{vec}(R\mathcal{J}(\hat{p}))$$

\rightsquigarrow a nonlinear least squares problem for R only.

- **Alternating projections (AP):**
alternate between solving two least squares problems

VARPRO is globally convergent with a super linear conv. rate.

AP is globally convergent with a linear convergence rate.

Structured low-rank approximation

No closed form solution is known for the general SLRA problem

$$\hat{p}^* := \arg \min_{\hat{p}} \|p - \hat{p}\| \quad \text{subject to} \quad \text{rank}(\mathcal{J}(\hat{p})) \leq r.$$

NP-hard, consider solution methods based on local optimization

Representing the constraint in a kernel form, the problem is

$$\min_{R, RR^\top = I_{m-r}} \left(\min_{\hat{p}} \|p - \hat{p}\| \quad \text{subject to} \quad R\mathcal{J}(\hat{p}) = 0 \right)$$

Note: Double minimization with bilinear equality constraint.

There is a matrix $G(R)$, such that $R\mathcal{J}(\hat{p}) = 0 \iff G(R)\hat{p} = 0$.

Software implementation

The structure of \mathcal{J} can be exploited for **efficient** $O(\dim(p))$ cost function and first derivative evaluations.

SLICOT library includes high quality FORTRAN implementation of algorithms for block Toeplitz matrices.

SLRA C software using I/O repr. and VARPRO approach
<http://www.esat.kuleuven.be/~imarkovs>

Based on the Levenberg–Marquardt alg. implemented in **MINPACK**.

Variations on low-rank approximation

• Cost functions

- weighted norms $(\text{vec}^\top(D)W\text{vec}(D))$
- information criteria $(\log \det(D))$

• Constraints and structures

- nonnegative
- sparse

• Data structures

- nonlinear models
- tensors

• Optimization algorithms

- convex relaxations

Nonnegative low-rank approximation

Constrained LRA arise in Markov chains and image mining

$$\min_{\hat{D}} \|D - \hat{D}\| \quad \text{subject to} \quad \text{rank}(\hat{D}) \leq r \text{ and } \hat{D}_{ij} \geq 0 \text{ for all } i, j.$$

Using an image representation, an **equivalent problem** is

$$\min_{P \in \mathbb{R}^{m \times r}, L \in \mathbb{R}^{r \times n}} \|D - PL\| \quad \text{subject to} \quad P_{ik}, L_{kj} \geq 0 \text{ for all } i, k, j.$$

Alternating projections algorithm:

- Choose an initial approximation $P^{(0)}$ and set $k := 0$.
- Solve: $L^{(k)} = \arg \min_L \|D - P^{(k)}L\|$ subject to $L \geq 0$.
- Solve: $P^{(k+1)} = \arg \min_P \|D - PL^{(k)}\|$ subject to $P \geq 0$.
- Repeat until convergence.

Weighted low-rank approximation

In the EIV model, LRA is ML assuming $\text{cov}(\text{vec}(\tilde{D})) = I$.

Motivation: incorporate prior knowledge W about $\text{cov}(\text{vec}(\tilde{D}))$

$$\min_{\hat{D}} \text{vec}^\top(D - \hat{D})W^{-1}\text{vec}(D - \hat{D}) \quad \text{subject to} \quad \text{rank}(\hat{D}) \leq r$$

Known in **chemometrics** as **maximum likelihood PCA**.

NP-hard problem, alternating projections is effective heuristic

Data fitting by a second order model

$$\mathcal{B}(A, b, c) := \{d \in \mathbb{R}^d \mid d^\top A d + b^\top d + c = 0\}, \quad \text{with } A = A^\top$$

Consider first **exact data**:

$$\begin{aligned} d \in \mathcal{B}(A, b, c) &\iff d^\top A d + b^\top d + c = 0 \\ &\iff \underbrace{\langle \text{col}(d \otimes_s d, d, 1), \text{col}(\text{vec}_s(A), b, c) \rangle}_{d_{\text{ext}} \quad \theta} = 0 \end{aligned}$$

$$\{d_1, \dots, d_N\} \in \mathcal{B}(\theta) \iff \theta \in \text{leftker} \underbrace{\begin{bmatrix} d_{\text{ext},1} & \dots & d_{\text{ext},N} \end{bmatrix}}_{D_{\text{ext}}}, \quad \theta \neq 0$$

$$\iff \text{rank}(D_{\text{ext}}) \leq d - 1$$

Therefore, for **measured data** \rightsquigarrow **LRA of D_{ext}** .

Notes:

- Special case \mathcal{B} an **ellipsoid** (for $A > 0$ and $4c < b^\top A^{-1}b$).
- Related to **kernel PCA**

Consistency in the errors-in-variables setting

Assume that the data is collected according to the EIV model

$$d_i = \bar{d}_i + \tilde{d}_i, \quad \text{where } \bar{d}_i \in \mathcal{B}(\bar{\theta}), \quad \tilde{d}_i \sim \mathcal{N}(0, \sigma^2 I).$$

LRA of D_{ext} (kernel PCA) \rightsquigarrow **inconsistent estimator**

$$\tilde{d}_{\text{ext},i} := \text{col}(\tilde{d}_i \otimes_s \tilde{d}_i, \tilde{d}_i, 0) \text{ is not Gaussian}$$

proposed method — incorporate bias correction in the LRA

Notes:

- works on the sample covariance matrix $D_{\text{ext}} D_{\text{ext}}^\top$
- the correction depends on the noise variance σ^2
- the core of the proposed method is the σ^2 estimator (possible link with methods for choosing **regularization par.**)

Rank minimization

Approximate modeling is a trade-off between:

- **fitting accuracy** and
- **model complexity**

Two possible scalarizations of the **bi-objective optimization** are:

LRA: minimize misfit under a constraint on complexity

RM: minimize complexity under a constraint (\mathcal{C}) on misfit

$$\min_X \text{rank}(X) \quad \text{subject to} \quad X \in \mathcal{C}$$

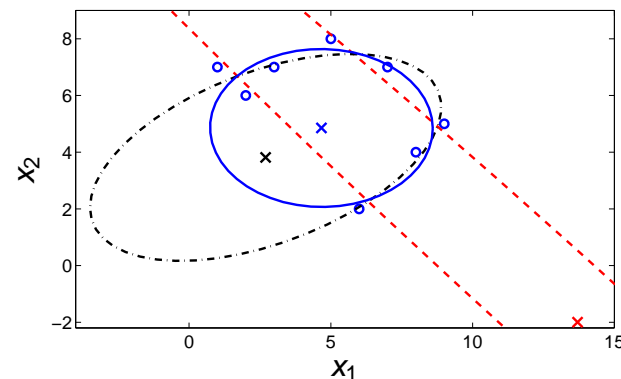
RM is also **NP-hard**, however, there are effective heuristics, e.g.,

with $X = \text{diag}(x)$, $\text{rank}(X) = \text{card}(x)$,

$$\ell_1 \text{ heuristic: } \min_x \|x\|_1 \quad \text{subject to} \quad \text{diag}(x) \in \mathcal{C}$$

Example: ellipsoid fitting

benchmark example of (Gander *et al.* 94), called “special data”



dashed — LRA solid — proposed method

dashed-dotted — orthogonal regression (geometric fitting)

○ — data points × — centers

Structured pseudospectra

$\Lambda(A)$ — the set of eigenvalues of $A \in \mathbb{C}^{n \times n}$

\mathbb{M} — a set of matrices ($\mathbb{M} = \{ \mathcal{S}(p) \mid p \in \mathbb{R}^{n_p} \}$)

Using the structured pseudospectra

$$\Lambda_\varepsilon(A) := \{ z \in \mathbb{C} \mid z \in \Lambda(B), B \in \mathbb{M}, \|A - B\|_2 \leq \varepsilon \}$$

one can determine the **distance to singularity**

$$d(A) := \min_{\Delta A \in \mathbb{M}} \|\Delta A\|_2 \quad \text{subject to} \quad A + \Delta A \text{ is singular}$$

which is a **special SLRA problem** with

1. square data matrix
2. perturbation measured by spectral norm, and
3. focus on minimum (vs minimizer) and singularity (vs rank).

Summary

- SLRA is a generic problem for data modeling.

search for more applications (pole placement, μ -analysis, ...)

- In general, SLRA is an NP-complete problem.

search for special cases that have “nice” solutions
e.g., circulant SLRA can be computed by DFT.

- The SLRA framework leads to conceptual unification.

Summary

- Efficient local solution methods

- Different rank representations (kernel, image, $AX = B$) lead to equivalent parameter optimization problems.

Computationally, however, these problems are different.

For example, the kernel representation leads to optimization on a Grassman manifold.

Currently, it is unexplored which parameterization is computational most beneficial.

Summary

- Effective heuristics, based on convex relaxations

- Practical advantage: one algorithm (and a piece of software) can solve a variety of problems

- Extensions of SLRA for tensors and nonlinear models

A framework with a potential for much to be done.

Thank you