

Chapter 8

Structured Low Rank Approximation

- Overview
- Low Rank Toeplitz Approximation
- Low Rank Circulant Approximation
- Low Rank Covariance Approximation
- Euclidean Distance Matrix Approximation
- Approximate GCD

Overview

- Given

- ◇ A target matrix $A \in R^{n \times n}$,
- ◇ An integer k , $1 \leq k < \text{rank}(A)$,
- ◇ A class of matrices Ω with linear structure,
- ◇ a fixed matrix norm $\| \cdot \|$;

Find

- ◇ A matrix $\hat{B} \in \Omega$ of rank k , and
- ◇

$$\|A - \hat{B}\| = \min_{B \in \Omega, \text{rank}(B)=k} \|A - B\|. \quad (1)$$

- Example of linear structure:

- ◇ Toeplitz or block Toeplitz matrices.
- ◇ Hankel or banded matrices.

- Applications:

- ◇ Signal and image processing with Toeplitz structure.
- ◇ Model reduction problem in speech encoding and filter design with Hankel structure.
- ◇ Regularization of ill-posed inverse problems.

Difficulties

- No easy way to characterize, either algebraically or analytically, a given class of structured lower rank matrices.
- Lack of explicit description of the feasible set \implies Difficult to apply classical optimization techniques.
- Little discussion on whether lower rank matrices with specified structure actually exist.

An Example of Existence

- Physics sometimes sheds additional light.
- The Toeplitz matrix

$$H := \begin{bmatrix} h_n & h_{n+1} & \dots & h_{2n-1} \\ \vdots & & & \vdots \\ h_2 & h_3 & \dots & h_{n+1} \\ h_1 & h_2 & \dots & h_n \end{bmatrix}$$

with

$$h_j := \sum_{i=1}^k \beta_i z_i^j, \quad j = 1, 2, \dots, 2n - 1,$$

where $\{\beta_i\}$ and $\{z_i\}$ are two sequences of arbitrary nonzero numbers satisfying $z_i \neq z_j$ whenever $i \neq j$ and $k \leq n$, is a Toeplitz matrix of rank k .

- The general Toeplitz structure preserving rank reduction problem as described in (1) remains open.
 - ◊ Existence of lower rank matrices of specified structure does not guarantee *closest* such matrices.
 - ◊ No $x > 0$ for which $1/x$ is minimum.

- For other types of structures, the existence question usually is a hard algebraic problem.

Another Hidden Catch

- The set of all $n \times n$ matrices with $\text{rank} \leq k$ is a *closed* set.
- The approximation problem

$$\min_{B \in \Omega, \text{rank}(B) \leq k} \|A - B\|$$

is *always* solvable, so long as the feasible set is non-empty.

- ◊ The rank condition is to be less than or equal to k , but not necessarily exactly equal to k .
- It is possible that a given target matrix A does not have a nearest rank k structured matrix approximation, but does have a nearest rank $k - 1$ or lower structured matrix approximation.

Low Rank Toeplitz Approximation

- Algebraic Structure of Low Rank Toeplitz Matrices.
- Constructing Low Rank Toeplitz Matrices.
 - ◇ Lift and Project Method
 - ◇ Parameterization by SVD
- Implicit Optimization
 - ◇ Engineerers' Misconception
 - ◇ Simplex Search Method
- Explicit Optimization
 - ◇ **constr** in MATLAB
 - ◇ **LANCELOT** on NEOS

A General Remark

- Introduce two procedures to tackle the structure preserving rank reduction problem numerically.
- The procedures can be applied to problems of any norm, any linear structure, and any matrix norm.
- Use the symmetric Toeplitz structure with Frobenius matrix norm to illustrate the ideas.

Algebraic Structure

- Identify a *symmetric* Toeplitz matrix by its first row,

$$T = T([t_1, \dots, t_n]) = \begin{bmatrix} t_1 & t_2 & \dots & t_n \\ t_2 & t_1 & \dots & t_{n-1} \\ \vdots & \dots & \dots & \\ t_{n-1} & & & t_2 \\ t_n & t_{n-1} & \dots & t_2 & t_1 \end{bmatrix}.$$

◇ \mathcal{T} = The affine subspace of all $n \times n$ symmetric Toeplitz matrices.

- Spectral decomposition of symmetric rank k matrices:

$$M = \sum_{i=1}^k \alpha_i \mathbf{y}^{(i)} \mathbf{y}^{(i)T}. \quad (2)$$

- Write $T = T([t_1, \dots, t_n])$ in terms of (2) \implies

$$\sum_{i=1}^k \alpha_i \mathbf{y}_j^{(i)} \mathbf{y}_{j+s}^{(i)} = t_{s+1}, \quad s = 0, 1, \dots, n-2, \quad 1 \leq j \leq n-s \quad (3)$$

◇ Lower rank matrices form an *algebraic variety*, i.e., solutions of polynomial systems.

Some Examples

- The case $k = 1$ is trivial.
 - ◊ Rank-one Toeplitz matrices form two simple one-parameter families,

$$T = \alpha_1 T([1, \dots, 1]), \quad \text{or}$$

$$T = \alpha_1 T([1, -1, 1, \dots, (-1)^{n-1}])$$

with arbitrary $\alpha_1 \neq 0$.

- For 4×4 symmetric Toeplitz matrices of rank 2, there are 10 unknowns in 6 equations.

$$\left\{ \begin{array}{l} \alpha_1 := \frac{\alpha_2 (y_1^{(2)2} - y_2^{(2)2})}{-y_1^{(1)2} + y_2^{(1)2}}, \\ y_3^{(1)} := \frac{y_2^{(1)} y_1^{(2)} y_1^{(1)} + 2 y_2^{(2)} y_2^{(1)2} - y_2^{(2)} y_1^{(1)2}}{y_2^{(1)} y_1^{(2)} + y_1^{(1)} y_2^{(2)}}, \\ y_4^{(1)} := \frac{-y_2^{(1)3} y_1^{(2)2} - 4 y_2^{(1)3} y_2^{(2)2} - 4 y_1^{(1)} y_1^{(2)} y_2^{(2)} y_2^{(1)2} - 2 y_2^{(1)} y_1^{(1)2} y_1^{(2)2} + 3 y_2^{(1)} y_2^{(2)2} y_1^{(1)2} + 2 y_1^{(2)} y_2^{(2)} y_1^{(1)3}}{y_2^{(1)2} y_1^{(2)2} + 2 y_2^{(1)} y_1^{(2)} y_1^{(1)} y_2^{(2)} + y_1^{(1)2} y_2^{(2)2}}, \\ y_3^{(2)} := \frac{-y_2^{(1)} y_1^{(2)2} - 2 y_2^{(1)} y_2^{(2)2} - y_1^{(2)} y_2^{(2)} y_1^{(1)}}{y_2^{(1)} y_1^{(2)} + y_1^{(1)} y_2^{(2)}}, \\ y_4^{(2)} := \frac{-3 y_2^{(1)2} y_1^{(2)2} y_2^{(2)} - 4 y_2^{(1)2} y_2^{(2)3} + 2 y_2^{(1)} y_1^{(1)} y_1^{(2)3} - 4 y_2^{(1)} y_1^{(1)} y_2^{(2)2} y_1^{(2)} - 2 y_2^{(2)} y_1^{(1)2} y_1^{(2)2} + y_1^{(1)2} y_2^{(2)3}}{y_2^{(1)2} y_1^{(2)2} + 2 y_2^{(1)} y_1^{(2)} y_1^{(1)} y_2^{(2)} + y_1^{(1)2} y_2^{(2)2}}. \end{array} \right.$$

- ◊ Explicit description of algebraic equations for higher dimensional lower rank symmetric Toeplitz matrices becomes unbearably complicated.

Let's See It!

- Rank deficient $T([t_1, t_2, t_3])$
 - ◇ $\det(T) = (t_1 - t_3)(t_1^2 + t_1t_3 - 2t_2^2) = 0$.
 - ◇ A union of two algebraic varieties.

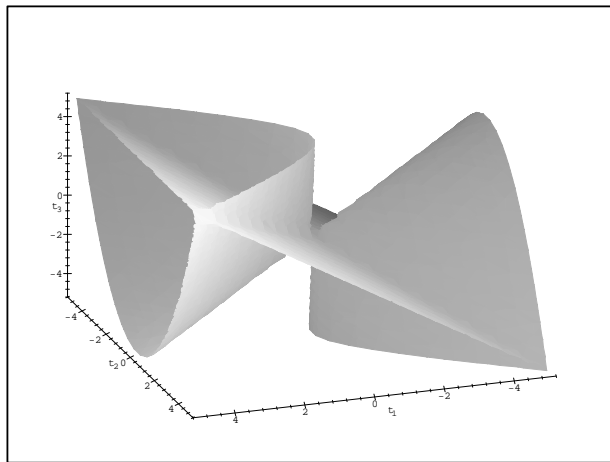


Figure 1: Lower rank, symmetric, Toeplitz matrices of dimension 3 identified in R^3 .

- The number of *local* solutions to the structured lower rank approximation problem is not unique.

Constructing Lower Rank Toeplitz Matrices

- Idea:
 - ◇ Rank k matrices in $R^{n \times n}$ form a *surface* $\mathcal{R}(k)$.
 - ◇ Rank k Toeplitz matrices = $\mathcal{R}(k) \cap \mathcal{T}$.
- Two approaches:
 - ◇ Parameterization by SVD:
 - ▷ Identify $M \in \mathcal{R}(k)$ by the triplet (U, Σ, V) of its singular value decomposition $M = U\Sigma V^T$.
 - U and V are orthogonal matrices, and
 - $\Sigma = \text{diag}\{s_1, \dots, s_k, 0, \dots, 0\}$ with $s_1 \geq \dots \geq s_k > 0$.
 - ▷ Enforce the structure.
 - ◇ Alternate projections between $\mathcal{R}(k)$ and \mathcal{T} to find intersections. (Cheney & Goldstein'59, Catzow'88)

Lift and Project Algorithm

- Given $A^{(0)} = A$, repeat projections until convergence:

- ◇ **LIFT**. Compute $B^{(\nu)} \in \mathcal{R}(k)$ nearest to $A^{(\nu)}$:

- ▷ From $A^{(\nu)} \in \mathcal{T}$, first compute its SVD

$$A^{(\nu)} = U^{(\nu)} \Sigma^{(\nu)} V^{(\nu)T}.$$

- ▷ Replace $\Sigma^{(\nu)}$ by $\text{diag}\{s_1^{(\nu)}, \dots, s_k^{(\nu)}, 0, \dots, 0\}$ and define

$$B^{(\nu)} := U^{(\nu)} \Sigma^{(\nu)} V^{(\nu)T}.$$

- ◇ **PROJECT**. Compute $A^{(\nu+1)} \in \mathcal{T}$ nearest to $B^{(\nu)}$:

- ▷ From $B^{(\nu)}$, choose $A^{(\nu+1)}$ to be the matrix formed by replacing the diagonals of $B^{(\nu)}$ by the averages of their entries.

- The general approach remains applicable to any other linear structure, and symmetry can be enforced.

- ◇ The only thing that needs to be modified is the projection in the projection (second) step.

Geometric Sketch

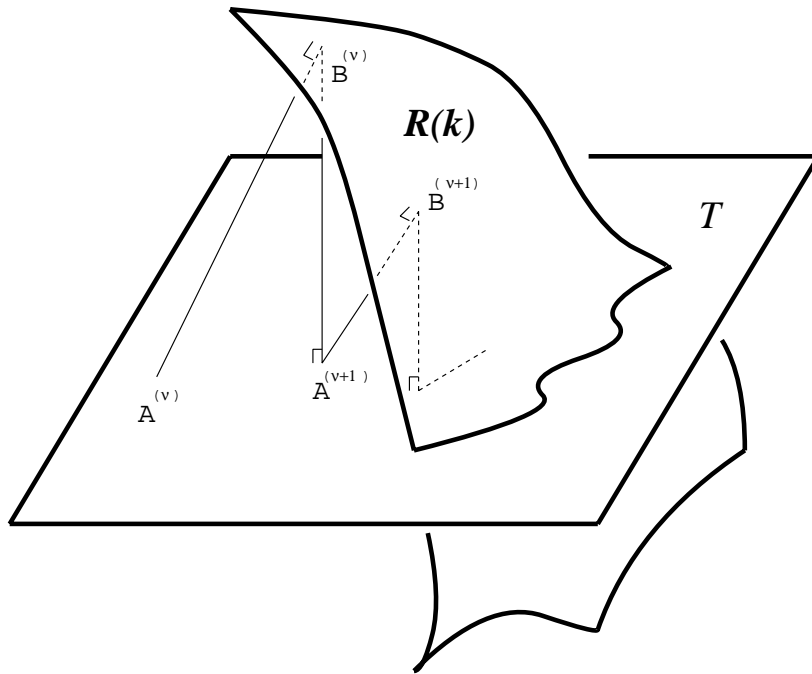


Figure 2: Algorithm 1 with intersection of lower rank matrices and Toeplitz matrices

Black-box Function

- Descent property:

$$\|A^{(\nu+1)} - B^{(\nu+1)}\|_F \leq \|A^{(\nu+1)} - B^{(\nu)}\|_F \leq \|A^{(\nu)} - B^{(\nu)}\|_F.$$

- ◇ Descent with respect to the Frobenius norm which is not necessarily the norm used in the structure preserving rank reduction problem.
- If all $A^{(\nu)}$ are distinct then the iteration converges to a Toeplitz matrix of rank k .
 - ◇ In principle, the iteration could be trapped in an impasse where $A^{(\nu)}$ and $B^{(\nu)}$ would not improve any more, but not experienced in practice.
- The lift and project iteration provides a means to define a *black-box function*

$$P : \mathcal{T} \longrightarrow \mathcal{T} \cap \mathcal{R}(k).$$

- ◇ The $P(\mathcal{T})$ is *presumably* piecewise continuous since all projections are continuous.

The graph of $P(T)$

- Consider $P : R^2 \longrightarrow R^2$:
 - ◇ Use the xy -plane to represent the domain of P for 2×2 symmetric Toeplitz matrices $T(t_1, t_2)$.
 - ◇ Use the z -axis to represent the image $p_{11}(T)$ and $p_{12}(T)$, respectively.

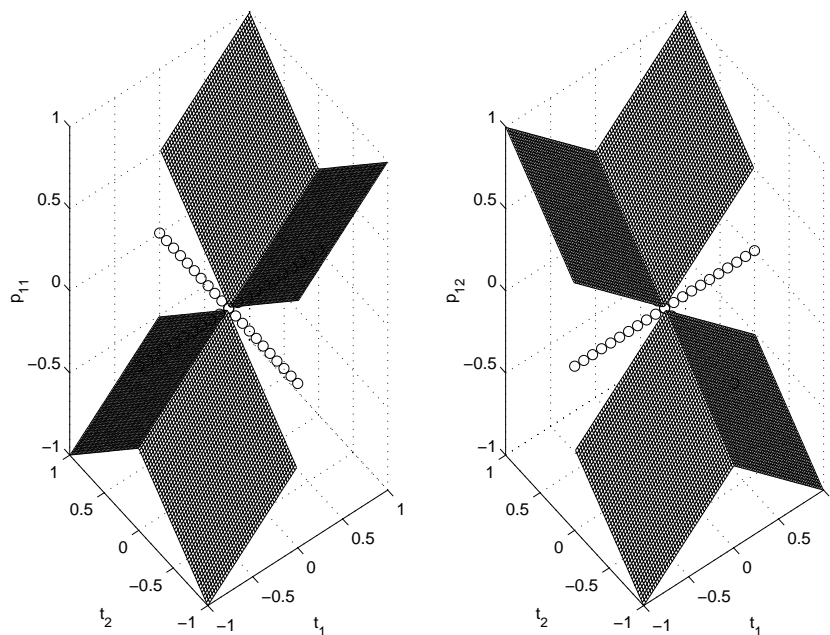


Figure 3: Graph of $P(T)$ for 2-dimensional symmetric Toeplitz T .

- Toeplitz matrices of the form $T(t_1, 0)$ or $T(0, t_2)$, corresponding to points on axes, converge to the zero matrix.

Implicit Optimization

- Implicit formulation:

$$\min_{T=\text{toeplitz}(t_1, \dots, t_n)} \|T_0 - P(T)\|. \quad (4)$$

- ◇ T_0 is the given target matrix.
 - ◇ $P(T)$, regarded as a black box function evaluation, provides a handle to manipulate the objective function $f(T) := \|T_0 - P(T)\|$.
 - ◇ The norm used in (4) can be any matrix norm.
- Engineers' misconception:
 - ◇ $P(T)$ is *not* necessarily the closest rank k Toeplitz matrix to T .
 - ◇ In practice, $P(T_0)$ has been used “as a cleansing process whereby any corrupting noise, measurement distortion or theoretical mismatch present in the given data set (namely, T_0) is removed.”
 - ◇ More needs to be done in order to find the *closest* lower rank Toeplitz approximation to the given T_0 as $P(T_0)$ is merely known to be in the feasible set.

Numerical Experiment

- An ad hoc optimization technique:
 - ◇ The simplex search method by Nelder and Mead requires only function evaluations.
 - ◇ Routine **fmins** in MATLAB, employing the simplex search method, is ready for use in our application.
- An example:
 - ◇ Suppose $T_0 = T(1, 2, 3, 4, 5, 6)$.
 - ◇ Start with $T^{(0)} = T_0$, and set worst case precision to 10^{-6} .
 - ◇ Able to calculate *all* lower rank matrices while maintaining the symmetric Toeplitz structure. Always so?
 - ◇ Nearly machine-zero of smallest calculated singular value(s) $\implies T_k^*$ is computationally of rank k .
 - ◇ T_k^* is only a local solution.
 - ◇ $\|T_k^* - T_0\| < \|P(T_0) - T_0\|$ which, though represents only a slight improvement, clearly indicates that $P(T_0)$ alone does not give rise to an optimal solution.

rank k	5	4	3	2	1
# of iterations	110	81	46	36	17
# of SVD calls	1881	4782	2585	2294	558
optimal solution	$\begin{bmatrix} 1.1046 \\ 1.8880 \\ 3.1045 \\ 3.9106 \\ 5.0635 \\ 5.9697 \end{bmatrix}$	$\begin{bmatrix} 1.2408 \\ 1.8030 \\ 3.0352 \\ 4.1132 \\ 4.8553 \\ 6.0759 \end{bmatrix}$	$\begin{bmatrix} 1.4128 \\ 1.7980 \\ 2.8171 \\ 4.1089 \\ 5.2156 \\ 5.7450 \end{bmatrix}$	$\begin{bmatrix} 1.9591 \\ 2.1059 \\ 2.5683 \\ 3.4157 \\ 4.7749 \\ 6.8497 \end{bmatrix}$	$\begin{bmatrix} 2.9444 \\ 2.9444 \\ 2.9444 \\ 2.9444 \\ 2.9444 \\ 2.9444 \end{bmatrix}$
$\ T_0 - T_k^*\ $	0.5868	0.9851	1.4440	3.2890	8.5959
singular values	$\begin{bmatrix} 17.9851 \\ 7.4557 \\ 2.2866 \\ 0.9989 \\ 0.6164 \\ 3.4638e-15 \end{bmatrix}$	$\begin{bmatrix} 17.9980 \\ 7.4321 \\ 2.2836 \\ 0.8376 \\ 2.2454e-14 \\ 2.0130e-14 \end{bmatrix}$	$\begin{bmatrix} 18.0125 \\ 7.4135 \\ 2.1222 \\ 1.9865e-14 \\ 9.0753e-15 \\ 6.5255e-15 \end{bmatrix}$	$\begin{bmatrix} 18.2486 \\ 6.4939 \\ 2.0884e-14 \\ 7.5607e-15 \\ 3.8479e-15 \\ 2.5896e-15 \end{bmatrix}$	$\begin{bmatrix} 17.6667 \\ 2.0828e-14 \\ 9.8954e-15 \\ 6.0286e-15 \\ 2.6494e-15 \\ 2.1171e-15 \end{bmatrix}$

Table 1: Test results for a case of $n = 6$ symmetric Toeplitz structure

Explicit Optimization

- Difficult to compute the gradient of $P(T)$.
- Other ways to parameterize structured lower rank matrices:
 - ◇ Use eigenvalues and eigenvectors for symmetric matrices;
 - ◇ Use singular values and singular vectors for general matrices.
 - ◇ Robust, but might have *overdetermined* the problem.

An Illustration

- Define

$$M(\alpha_1, \dots, \alpha_k, \mathbf{y}^{(1)}, \dots, \mathbf{y}^{(k)}) := \sum_{i=1}^k \alpha_i \mathbf{y}^{(i)} \mathbf{y}^{(i)T}.$$

- Reformulate the symmetric Toeplitz structure preserving rank reduction problem *explicitly* as

$$\begin{aligned} \min \quad & \|T_0 - M(\alpha_1, \dots, \alpha_k, \mathbf{y}^{(1)}, \dots, \mathbf{y}^{(k)})\| \quad (5) \\ \text{subject to} \quad & m_{j,j+s-1} = m_{1,s}, \quad (6) \\ & s = 1, \dots, n-1, \\ & j = 2, \dots, n-s+1, \end{aligned}$$

if $M = [m_{ij}]$.

- ◇ Objective function in (5) is described in terms of the non-zero eigenvalues $\alpha_1, \dots, \alpha_k$ and the corresponding eigenvectors $\mathbf{y}^{(1)}, \dots, \mathbf{y}^{(k)}$ of M .
 - ◇ Constraints in (6) are used to ensure that M is symmetric and Toeplitz.
- For other types of structures, we only need modify the constraint statement accordingly.
- The norm used in (5) can be arbitrary but is fixed.

Redundant Constraints

- Symmetric centro-symmetric matrices have special spectral properties:
 - ◇ $\lceil n/2 \rceil$ of the eigenvectors are symmetric; and
 - ◇ $\lfloor n/2 \rfloor$ are skew-symmetric.
 - ▷ $v = [v_i] \in \mathbb{R}^n$ is symmetric (or skew-symmetric) if $v_i = v_{n-i}$ (or $v_i = -v_{n-i}$).
- Symmetric Toeplitz matrices are symmetric and centro-symmetric.
- The formulation in (5) does not take this spectral structure into account in the eigenvectors $y^{(i)}$.
 - ◇ More variables than needed have been introduced.
 - ◇ May have overlooked any internal relationship among the $\frac{n(n-1)}{2}$ equality constraints.
 - ◇ May have caused, inadvertently, additional computation complexity.

Using Existing Optimization Codes

- Using **constr** in MATLAB

- ◇ Routine **constr** in MATLAB:

- ▷ Uses a sequential quadratic programming method.
- ▷ Solve the Kuhn-Tucker equations by a quasi-Newton updating procedure.
- ▷ Can estimate derivative information by finite difference approximations.
- ▷ Readily available in Optimization Toolbox.

- ◇ Our experiments:

- ▷ Use the same data as in the implicit formulation.
- ▷ Case $k = 5$ is computationally the same as before.
- ▷ Have trouble in cases $k = 4$ or $k = 3$,
 - Iterations will not improve approximations at all.
 - MATLAB reports that the optimization is terminated successfully.

- Using **LANCELOT** on NEOS

- ◇ Reasons of failure of MATLAB are not clear.
 - ▷ Constraints might no longer be linearly independent.
 - ▷ Termination criteria in **constr** might not be adequate.
 - ▷ Difficult geometry means hard-to-satisfy constraints.
- ◇ Using more sophisticated optimization packages, such as **LANCELOT**.
 - ▷ A standard Fortran 77 package for solving large-scale nonlinearly constrained optimization problems.
 - ▷ Break down the functions into sums of *element functions* to introduce sparse Hessian matrix.
 - ▷ Huge code. See

<http://www.rl.ac.uk/departments/ccd/numerical/lancelot/sif/sifhtml.html>.

- ▷ Available on the NEOS Server through a socket-based interface.
- ▷ Uses the **ADIFOR** automatic differentiation tool.

◇ **LANCELOT** works.

- ▷ Find optimal solutions of problem (5) for all values of k .
- ▷ Results from **LANCELOT** agree, up to the required accuracy 10^{-6} , with those from **fmins**.
- ▷ Rank affects the computational cost nonlinearly.

rank k	5	4	3	2	1
# of variables	35	28	21	14	7
# of f/c calls	108	56	47	43	19
total time	12.99	4.850	3.120	1.280	.4300

Table 3: Cost overhead in using **LANCELOT** for $n = 6$.

Conclusions

- Structure preserving rank reduction problems arise in many important applications, particularly in the broad areas of signal and image processing.
- Constructing the nearest approximation of a given matrix by one with any rank and any linear structure is difficult in general.
- We have proposed two ways to formulate the problems as standard optimization computations.
- It is now possible to tackle the problems numerically via utilizing standard optimization packages.
- The ideas were illustrated by considering Toeplitz structure with Frobenius norm.
- Our approach can be readily generalized to consider rank reduction problems for any given linear structure and of any given matrix norm.

Low Rank Circulant Approximation

- Basic Properties
- (Inverse) Eigenvalue Problem
- Conjugate Evenness
- Low Rank Approximation
- Tree Structure
- Numerical Experiment

Basic Properties

- A circulant matrix $C = \text{Circul}(c)$

$$C = \begin{bmatrix} c_0 & c_1 & \dots & c_{n-1} \\ c_{n-1} & c_0 & c_1 & \dots & c_{n-2} \\ c_{n-2} & c_{n-1} & c_0 & \dots & c_{n-3} \\ \vdots & \vdots & \dots & \dots & \vdots \\ c_1 & c_2 & \dots & c_{n-1} & c_0 \end{bmatrix}$$

is uniquely determined by the first row c .

◇ Define

$$\Pi := \begin{bmatrix} 0 & 1 & 0 & \dots & 0 \\ 0 & 0 & 1 & & 0 \\ \vdots & & \dots & \dots & \vdots \\ 0 & & & & 1 \\ 1 & 0 & & \dots & 0 \end{bmatrix}.$$

Then

$$\text{Circul}(c) = \sum_{k=0}^{n-1} c_k \Pi^k = P_c(\Pi)$$

with characteristic polynomial

$$P_c(x) = \sum_{k=0}^{n-1} c_k x^k.$$

Elementary Spectral Properties

- Define

$$\Omega := \text{diag}(1, \omega, \omega^2, \dots, \omega^{n-1}), \quad \omega := \exp\left(\frac{2\pi i}{n}\right).$$

- Define the Fourier matrix F where

$$F^* := \frac{1}{\sqrt{n}} \begin{bmatrix} 1 & 1 & 1 & \dots & 1 \\ 1 & \omega & \omega^2 & \dots & \omega^{n-1} \\ 1 & \omega^2 & \omega^4 & \dots & \omega^{2n-2} \\ \vdots & & & & \vdots \\ 1 & \omega^{n-1} & \omega^{n-2} & \dots & \omega \end{bmatrix}.$$

- ◊ F is unitary.

- The forward shift matrix Π is unitarily diagonalizable.

$$\Pi = F^* \Omega F.$$

- The circulant matrix $Circul(c)$ with any given row vector c has a spectral decomposition

$$Circul(c) = F^* P_c(\Omega) F.$$

- Spectral properties:
 - ◇ Closely related to the discrete Fourier transform.
 - ◇ Explicit solution for eigenvalue and inverse eigenvalue problems.
 - ◇ FFT calculation.

(Inverse) Eigenvalue Problem

- Forward problem:

- ◇ Eigenvalues of $Circul(c)$:

$$\lambda = [P_c(1), \dots, P_c(\omega^{n-1})].$$

- ◇ Can be computed from

$$\lambda^T = \sqrt{n} F^* c^T.$$

- Inverse problem:

- ◇ Given any vector $\lambda := [\lambda_0, \dots, \lambda_{n-1}] \in C^n$, define

$$c^T = \frac{1}{\sqrt{n}} F \lambda^T.$$

- ◇ $Circul(c)$ has eigenvalues in vector λ .

- Both matrix-vector products involved done in $O(n \log_2 n)$ flops.
- If all the eigenvalues are distinct, there are precisely $n!$ distinct circulant matrices with the prescribed spectrum.

Conjugate Evenness

- $c^T = \frac{1}{\sqrt{n}}F\lambda^T$ is real if and only if $\lambda^T = \sqrt{n}F^*c^T$ is conjugate-even.

- ◇ If $n = 2m$,

$$\lambda = [\lambda_0, \lambda_1, \dots, \lambda_{m-1}, \lambda_m, \overline{\lambda_{m-1}}, \dots, \overline{\lambda_1}].$$

- ▷ $\lambda_0, \lambda_m \in \mathbb{R}$. (Absolutely real, others real or complex.)

- ◇ If $n = 2m + 1$,

$$\lambda := [\lambda_0, \lambda_1, \dots, \lambda_m, \overline{\lambda_m}, \dots, \overline{\lambda_1}].$$

- ▷ $\lambda_0 \in \mathbb{R}$. (Absolutely real.)

- Singular value decomposition of $Circul(c)$:

$$Circul(c) = (F^*P_c(\Omega)|P_c(\Omega)|^{-1})|P_c(\Omega)|F$$

- ◇ Singular values are $|P_c(\omega^k)|, k = 0, 1, \dots, n - 1$.

- ◇ At most $\lceil \frac{n+1}{2} \rceil$ distinct singular values.

Low Rank Approximation

- Given $A \in R^{n \times n}$, its nearest circulant matrix approximation $Circul(c)$ is given by the projection (T. Chan)

$$c_k := \frac{1}{n} \langle A, \Pi^k \rangle, \quad k = 0, \dots, n-1,$$

- ◇ $Circul(c)$ is generally of **full rank** even if A has lower rank.
- How to reduce the rank of $Circul(c)$?
 - ◇ The truncated singular value decomposition (TSVD) gives rise to the nearest low rank approximation in Frobenius norm.
 - ◇ The TSVD of $Circul(c)$ is automatically circulant.
 - ◇ But: the TSVD can lead to a **complex** circulant approximation.

Trivial $O(n \log n)$ TSVD Algorithm

- Given a real matrix A and a fixed rank $\ell \leq n$,
 1. Use the projection to find the nearest real circulant matrix approximation $Circul(c)$ of A .
 2. Use the FFT to calculate the spectrum λ of the matrix $Circul(c)$.
 3. Arrange all elements of $|\lambda|$ in descending order, including those with equal modulus.
 4. Let $\hat{\lambda}$ be the vector consisting of elements of λ , but those corresponding to the last $n - \ell$ singular values in the descending order are set to zero.
 5. Apply the inverse FFT to $\hat{\lambda}$ to determine a nearest circulant matrix $Circul(\hat{c})$ of rank ℓ to A .
- The resulting matrix $Circul(\hat{c})$ is complex-valued in general.
 - ◇ Need to preserve the conjugate-even structure.
 - ◇ Need to modify the TSVD strategy.

Data Matching Problem

- The low rank “real” circulant approximation problem is equivalent to a data matching problem:

(DMP) *Given a conjugate-even vector $\lambda \in C^n$, find its nearest conjugate-even approximation $\hat{\lambda} \in C^n$ subject to the constraint that $\hat{\lambda}$ has exactly $n - \ell$ zeros.*

- How to solve the DMP?

- ◊ Write $\hat{\lambda} = [\hat{\lambda}_1, 0] \in C^n$ with $\hat{\lambda}_1 \in C^\ell$ being arbitrary.

- ◊ Consider the problem of minimizing

$$F(P, \hat{\lambda}) = \|P\hat{\lambda}^T - \lambda^T\|^2$$

with a permutation matrix P .

- ▷ P is used to search for the match.

- ◊ Write $P = [P_1, P_2]$ with $P_1 \in R^{n \times \ell}$.

◇ A least squares problem:

$$F(P, \hat{\lambda}) = \|P_1 \hat{\lambda}_1^T - \lambda^T\|^2$$

◇ The optimal solution is

$$\hat{\lambda}_1 = \lambda P_1.$$

▷ The entries of $\hat{\lambda}_1$ must be a portion of λ .

◇ The objective function becomes

$$F(P, \hat{\lambda}) = \|(P_1 P_1^T - I)\lambda\|^2.$$

▷ $P_1 P_1^T - I$ is just a projection.

▷ The optimal permutation P should be such that $P_1 P_1^T$ projects λ to its first ℓ components with largest modulus.

- Without the conjugate-even constraint, the answer to the data matching problem corresponds precisely to the usual TSVD selection criterion.
- With the conjugate-even constraint, the above criterion remains effective subject to the conjugate-even structure inside λ .

An Example using a Tree Structure

- Consider the case $n = 6$.
- Suppose λ_1, λ_2 are complex and distinct.
- Six possible conjugate-even structures.
- Tree graph:
 - ◇ Each node in the tree represents an element of λ .
 - ◇ Arrange the nodes from top to bottom in descending order of their moduli.
 - ◇ In case of a tie,
 - ▷ Complex conjugate nodes stay at the same level.
 - ▷ Real node is below the complex nodes.
- If $\lambda_1, \overline{\lambda_1}, \lambda_0, \lambda_2, \overline{\lambda_2}, \lambda_3$, then the tree is given by:

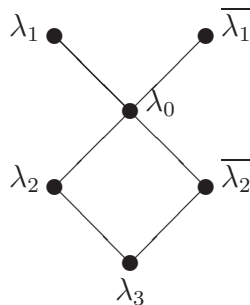


Figure 4: Tree graph of $\lambda_1, \overline{\lambda_1}, \lambda_0, \lambda_2, \overline{\lambda_2}, \lambda_3$ with $|\lambda_1| \geq |\lambda_0| > |\lambda_2| \geq |\lambda_3|$.

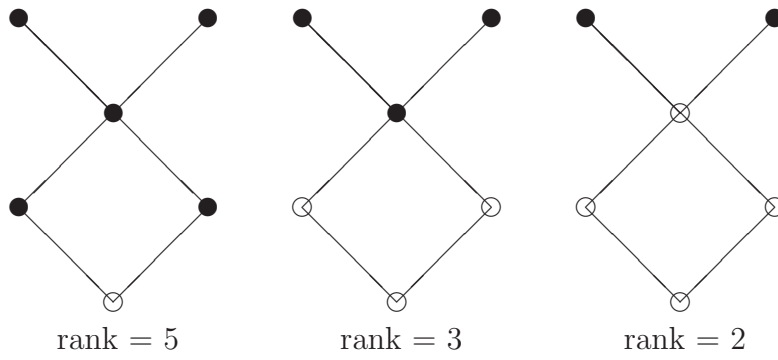


Figure 5: Tree graphs of $\hat{\lambda}$ with rank 5, 3, and 2.

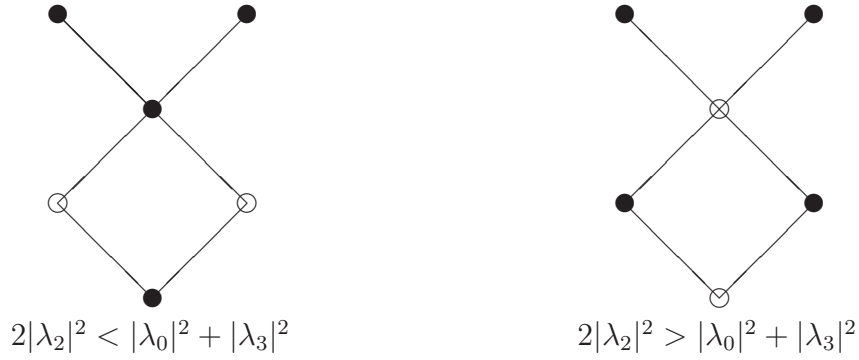


Figure 6: Tree graphs of $\hat{\lambda}$ with rank 4.

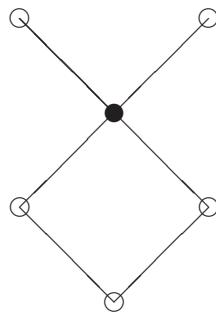


Figure 7: Tree graph of $\hat{\lambda}$ with rank 1.

rank λ	5	4	3	2	1	other possibilities	

Figure 8: Possible solutions to the DMP when $n = 6$.

Numerical Experiment

- The low rank approximation algorithm needs to be smart enough to explore the conjugate-even structure, to truncate, and to reassemble the conjugate-even structure.
- Numerical complexity is $O(n \log n)$ flops.

Example 1: Symmetric Illustration

Consider the 8×8 symmetric $Circul(c)$:

$$c = [0.5404, 0.2794, 0.1801, -0.0253, -0.2178, -0.0253, 0.1801, 0.2794].$$

- Eigenvalues (in descending order):

$$[1.1909, 1.1892, 1.1892, 0.3273, 0.3273, \mathbf{0.1746}, -0.0376, -0.0376]$$

- For rank 7 approximation, the usual TSVD would set -0.0376 to zero, resulting in a complex matrix.
- Use the conjugate-even eigenvalues

$$\hat{\lambda} = [1.1909, 1.1892, 0.3273, -0.0376, \mathbf{0} -0.0376, 0.3273, 1.1892],$$

to obtain the best real-valued, rank 7, approximation $Circul(\hat{c})$ via the FFT:

$$\hat{c} = [0.5186, 0.3657, 0.0670, -0.0680, -0.0572, -0.0680, 0.0670, 0.3657].$$

- To obtain the best real-value, rank 4, circulant approximation, use eigenvalues $\hat{\lambda}$

$$\hat{\lambda} = [1.1909, 1.1892, 0, 0, 0.3273, 0, 0, 1.1892].$$

Example 2: Complex Illustration

Consider the 9×9 $Circul(c)$ with

$$c = [1.6864, 1.7775, 1.9324, 2.9399, 1.9871, 1.7367, 4.0563, 1.2848, 2.5989].$$

- Eigenvalues: structure given by

$$[20.0000, \\ -2.8130 + 1.9106i, -2.8130 - 1.9106i, 3.0239 - 1.0554i, 3.0239 + 1.0554i, \\ -1.3997 + 0.7715i, -1.3997 - 0.7715i, -1.2223 - 0.2185i, -1.2223 + 0.2185i].$$

- To obtain a real-valued circulant approximation of rank 8, we have no choice but to select the set the *largest* eigenvalue (singular value) of $Circul(c)$ to zero.
 - ◇ Setting the largest singular value to zero to obtain the nearest real low rank approximation is quite counter-intuitive to the usual sense of TSVD.
 - ◇ Apply algorithm to reduce the rank further, to 7.

Example 3: Perturbed Case

- Let $C_\kappa \in R^{n \times n}$ be a given circulant matrix of rank κ . Random noise added to C_κ will destroy the circulant structure as well as the rank condition.
- Let $E \in R^{n \times n}$ denote a random but fixed circulant matrix with unit Frobenius norm, and let

$$W_j = C_\kappa + 10^{-j}E, \quad j = 1, \dots, 12.$$

- W_j will almost certainly be of full rank. Note that $\|W_j - C_\kappa\| = 10^{-j}$. It will be interesting to see if W_j has any closer circulant matrix approximation of rank κ other than C_κ , especially when j is large.
- Test case with $n = 100$, $\kappa = 73$, and a predetermined matrix C_{73} .
- Using our algorithm to find the best circulant approximation Z_j to W_j , we find that it is always the case that

$$\|W_j - Z_j\| < \|W_j - C_\kappa\|$$

for all j , i.e., our real circulant approximation is closest.

Conclusion

- For any given real data matrix, its nearest real circulant approximation can simply be determined from the average of its diagonals.
- The nearest low rank (possibly complex) approximation to the circulant matrix can be determined effectively from the TSVD and the FFT.
- To construct a real circulant matrix with specified spectrum, the eigenvalues must appear in conjugate-even form. So, the truncation criteria for a nearest low rank, real, circulant matrix approximation must be modified.
- We have proposed a fast algorithm with $O(n \log n)$ complexity to accomplish all of these objectives.

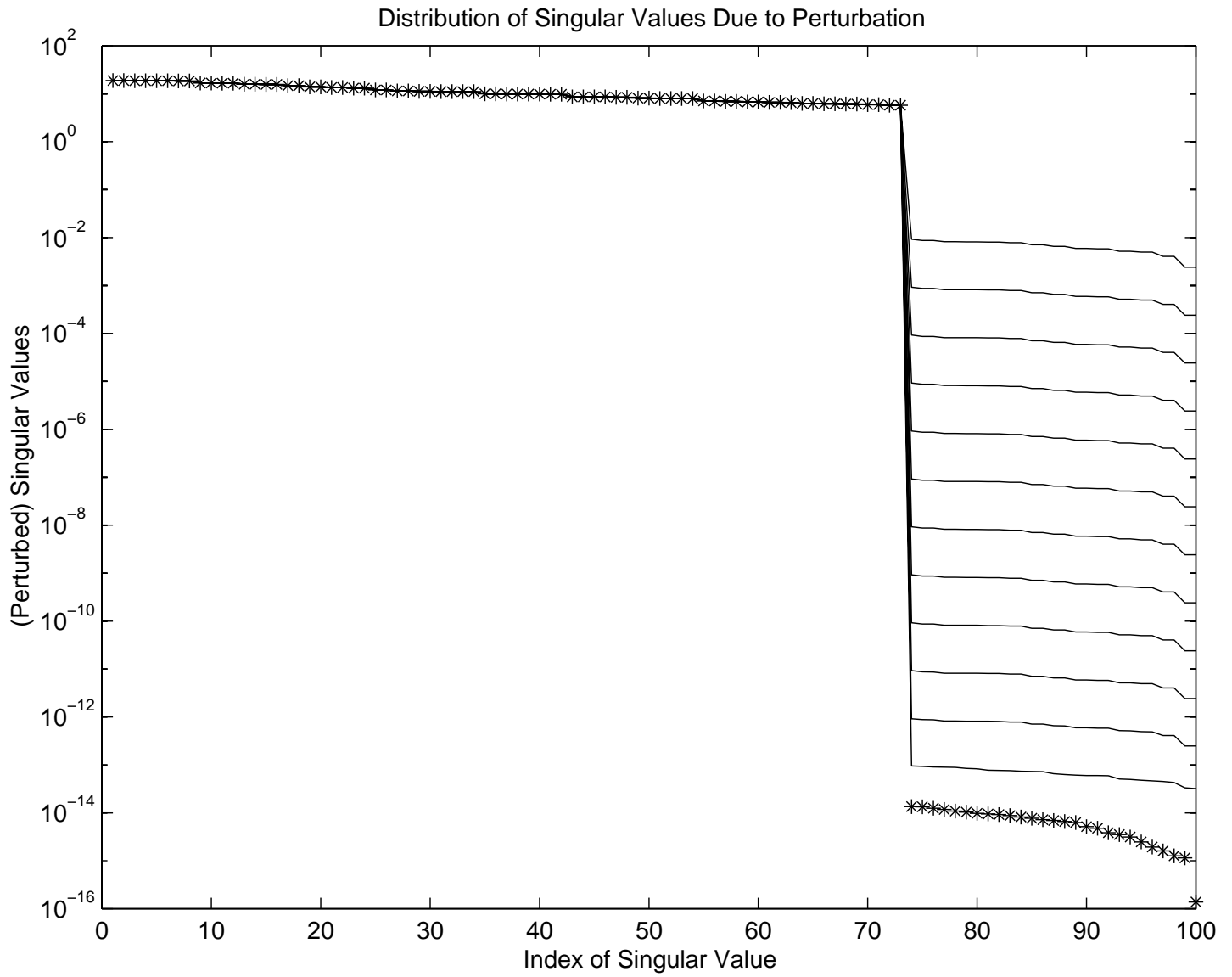


Figure 9: Distribution of Singular Values.

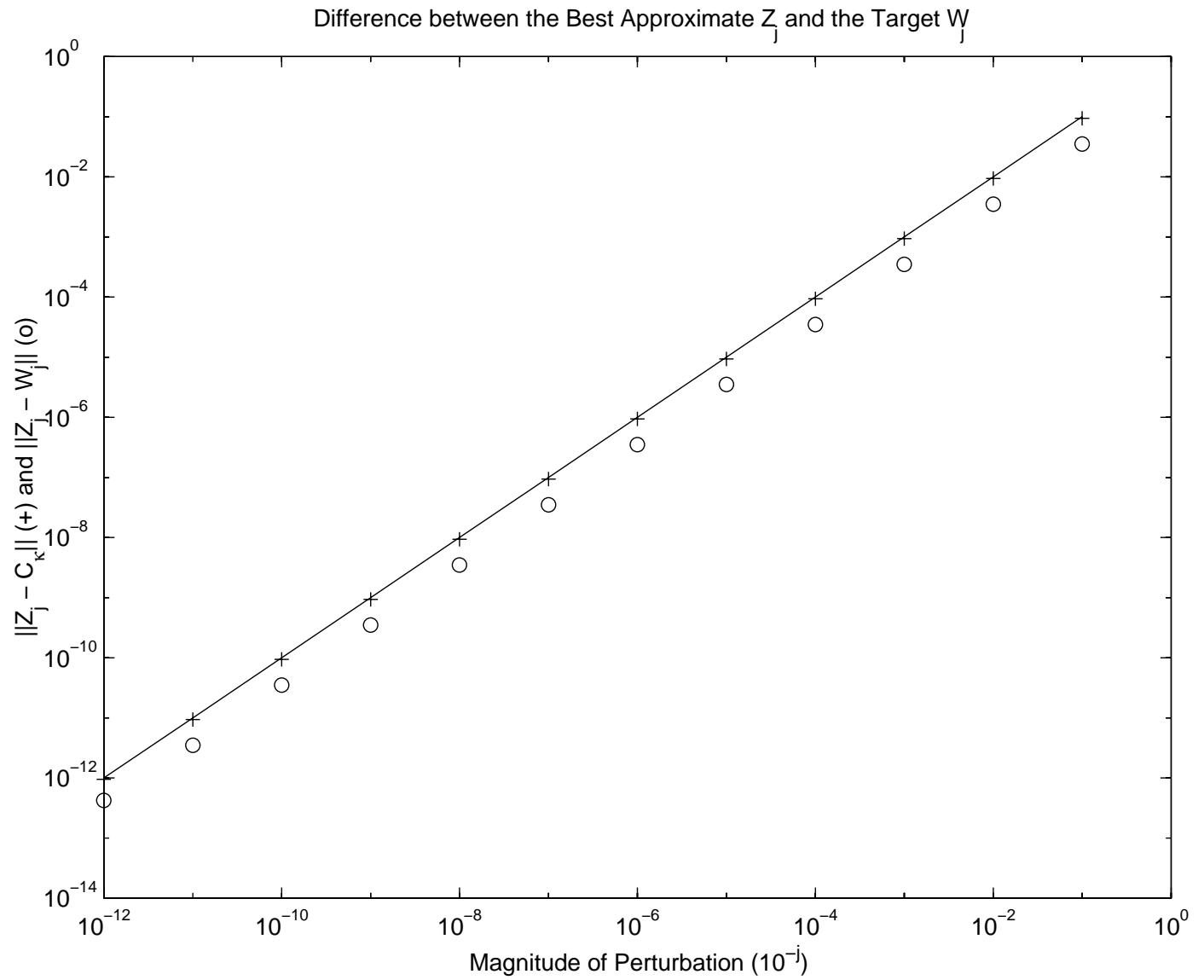


Figure 10: Errors in Approximation.

Lor Rank Covariance Approximation

Euclidian Distance Matrix Approximation

Approximate GCD

f-COUNT	FUNCTION	MAX{g}	STEP	Procedures
29	0.958964	8.65974e-15	1	
77	0.958964	2.66454e-14	1.91e-06	
131	0.958964	2.70894e-14	2.98e-08	Hessian modified twice
185	0.958964	2.70894e-14	2.98e-08	
239	0.958964	2.73115e-14	2.98e-08	
289	0.958964	2.77556e-14	4.77e-07	
337	0.958964	2.77556e-14	1.91e-06	
393	0.958964	2.77556e-14	7.45e-09	Hessian modified twice
445	0.958964	5.28466e-14	1.19e-07	
501	0.958964	5.68434e-14	7.45e-09	
557	0.958964	5.70655e-14	7.45e-09	Hessian not updated
613	0.958964	5.66214e-14	7.45e-09	
667	0.958964	5.55112e-14	2.98e-08	Hessian modified twice
713	0.958964	3.17302e-13	7.63e-06	
761	0.958964	2.61569e-13	1.91e-06	
812	0.958964	2.60014e-13	-2.38e-07	Hessian modified twice
856	0.958964	2.57794e-13	3.05e-05	Hessian modified twice
900	0.958964	2.56462e-13	3.05e-05	Hessian modified twice
948	0.958964	2.57128e-13	1.91e-06	
994	0.958964	2.56684e-13	7.63e-06	
1038	0.958964	3.42837e-13	3.05e-05	
1083	0.958964	3.41727e-13	-1.53e-05	Hessian modified twice
1124	0.958964	3.92575e-13	0.000244	Hessian modified twice
1161	0.958964	5.04485e-13	0.00391	Hessian modified twice
1200	0.958964	5.12923e-13	0.000977	Hessian modified twice
1233	0.958964	5.61551e-13	0.0625	Hessian modified twice
1272	0.958964	5.86642e-13	0.000977	Hessian modified twice
1308	0.958964	4.84279e-13	0.00781	Hessian modified twice
1309	0.958964	4.84723e-13	1	Hessian modified twice

Optimization Converged Successfully

Table 2: A typical output of intermediate results from **constr**.