

## Parallel Iterations for Cholesky and Rohn’s Methods for Solving over Determined Linear Interval Systems

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**Abstract:** The aim is to present Krawczyk’s iteration in the context of parallel Rohn’s method and another parallel Interval Cholesky method to improve results for the solution of over determined linear interval system. We compare note with results obtained from using the Block Krawczyk iterative method introduced by Rump, and presented in details by Popova, which forms the basis of discussions. It is shown that parallel Rohn’s method which is independent of interval data input gave narrowest interval width of the solution set to the over determined interval linear system followed by Interval Cholesky method while Block Krawczyk’s method given in Popova, has the largest interval width.

**Key words:** AMS Subject Classification (2000), interval Cholesky method, Krawczyk’s method, over determined linear interval system, Rohn’s method, 65G20 and 65G30

### INTRODUCTION

We signify with the following notation  $\mathbb{R}^n$ ,  $\mathbb{IR}^n$ ,  $\mathbb{R}^{n \times n}$ ,  $\mathbb{IR}^{n \times n}$  as representing the sets of real vectors, interval vectors, real matrices and interval matrices, respectively. We assume that the interval matrix  $A \in [A]$  is an  $m \times n$  matrix where  $m > n$ .

We are interested in the solution of over determined linear system with interval data:

$$Ax = b \tag{1}$$

where the interval entries of A and b are described by:

$$A = (A_c - \Delta, A_c + \Delta)$$

$$b = (b_c - \delta, b_c + \delta)$$

Such that:

$$A_c = \frac{1}{2} \left( \underline{A} + \bar{A} \right)$$

$$x_c = \frac{1}{2} \left( \underline{x} + \bar{x} \right)$$

$$b_c = \frac{1}{2} \left( \underline{b} + \bar{b} \right)$$

The terms  $\Delta$  and  $\delta$  denote respectively the radius an interval matrix and interval vector. Such situation arises when we describe a dynamical system from state

$$S(a) = (S_1(a), \dots, S_n(a))$$

at a given time to the state

$$S(a+1) = (S_1(a+1), \dots, S_n(a+1))$$

at the epoch of time a+1 which leads to dependence

$$S_n(a+1) = f_i (S_1(a), \dots, S_n(a))$$

This means that the parameters describing the coefficients of the linear system are afflicted with uncertainties.

We define the set:

$$\Sigma_{m \times n} = \{ x_c \in \mathbb{R}^n \mid A \in [A], \exists b \in [b],$$

$$x_c = \operatorname{argmin}_{x \in \mathbb{R}^n} \|b - Ax\| \}$$

$$= \{ x \in \mathbb{R}^n \mid A \in [A], \exists b \in [b], A^H Ax = A^H b \}$$

(Rump, 1993) proposed and solved the over determined system by considering the augmented square linear system:

$$\begin{pmatrix} [A] & -I \\ 0 & [A]^H \end{pmatrix} \begin{pmatrix} x \\ y \end{pmatrix} = \begin{pmatrix} [b] \\ 0 \end{pmatrix}, m > n \tag{2}$$

Following the approach presented in (Rump, 1993), the following theorem was proposed by (Popova, 2006).

**Theorem 1:** (Popova, 2006).

Let  $[A] \in \mathbb{IR}^{m \times n}, [b] \in \mathbb{IR}^n, m > n$

Defining:

$$[B] \in \mathbb{IR}^{((m+n) \times (m+n))}, [V] \in \mathbb{IR}^{m+n}$$

representing a square interval matrix, respectively, the interval vector in (2) and letting,

$$R \in \mathbb{R}^{(m+n) \times (m+n)}, \bar{U} \in \mathbb{R}^{m+n}$$

(i) Let  $[Z] \in \mathbb{IR}^{m+n}$  be defined in the form:

$$\begin{aligned}
 [Z]_{l=1}^{m+n} &= \sum_{\mu}^m r_{i\mu} (b_{i\mu} + \bar{\mu}_{n+\mu}) \\
 &- \sum_{w=l}^m \sum_{v=l}^n (r_{i\mu} \bar{U}_v + r_{i,m+\nu} \bar{U}_{n+\mu}) R_{\mathbb{E}}([a_{\mu v}]) \\
 &- \sum_{\mu=1}^m \sum_{v=1}^n (r_{i\mu} \bar{\mu}_v - r_{i,m+\nu} \bar{\mu}_{n-w}) I_m([a_{\mu v}])
 \end{aligned}$$

(ii) Let  $[C] \in \mathbb{IR}^{(m+n) \times (m+n)}$  be defined by I-R[B] and

$$\text{let } [u] \in \mathbb{IR}^{m+n}.$$

Assuming  $[w] \in \mathbb{IR}^{m+n}$ , where

$$[w_i]_{i=1}^{m+n} = \{ [Z] + [C] \cdot [uu] \}_{i,l}$$

$$[uu] = ([w_1], \dots, [w_{i-1}], [u_i], \dots, [u_{m+n}])^T$$

If  $[w] \subseteq_{\neq} [u]$  then every matrix  $A \in [A]$  has full rank  $n$ , and for every  $A \in [A]$ ,  $b \in [b]$  with  $x_c$  minimizing  $\|b - Ax\|$ , the unique solution  $x_c$  satisfies  $\hat{x} \in x_c + [x]$  where  $\hat{x}$  and  $[x]$  are the first  $n$  components of  $u_c$ , resp  $[w]$ .

(iii) With  $[\Delta] = [C][w] \in \mathbb{IR}^{m+n}$  and  $[xx] \in \mathbb{IR}^n$  defined by

$$[xx_i] = \inf([z_i]) + \text{Sup}([\Delta_i])$$

$$\sup([z_i] + \inf([\Delta_i])), 1 \leq i \leq n,$$

The following inner and outer estimations were obtained by (Popova, 2006) in the form:

$$x_c + [xx] \subseteq (\sum_{m \times n}) \subseteq x_c + [x]$$

In this work, we are interested in using parallel Cholesky Factorization for the algebraic inclusion of over determined linear interval systems of equations. A procedure for speeding up the Interval Cholesky

Factorization using a method described in (Gargantini, 1980) in conjunction with (Carstensen and Petkovic, 1994) and of course (Rump, 1999) to narrow the overestimation of resulting disks is implied.

We introduce a theorem due to (Rohn, 1996) which is free of interval input in section three which can be interpreted as a variant of (Rump, 1993) approach. The convergence of Rohn's method is dependent on the existence of an approximation matrix of the Moore penrose inverse of A.

Finally, numerical example taken from (Popova, 2006) is used to demonstrate the quality of inclusion bounds provided by these methods. We concluded the paper based on the findings with results from these methods.

## MATERIALS AND METHODS

We use parallel interval Cholesky method and parallel (Rohn, 1996)'s method as our basis of experiments. Below we give mathematical details of the methods.

**Parallel interval Cholesky method:** We discuss parallel interval Cholesky method by replacing real entries and arithmetic operations with interval ones. The idea is based on the extended interval arithmetic, which dates back to (Gargantini, 1976) and further studied by (Rump, 1993). Using the notation

$$[A] = \begin{bmatrix} A, \bar{A} \\ - \end{bmatrix} = a_{ij} = \begin{bmatrix} a, \bar{a}_{ij} \\ - \bar{a}_{ij} \end{bmatrix}$$

where, we can convert the over determined system (1) into equivalent interval linear system of size  $n \times n$  coefficient matrix by a certain matrix R called a preconditioner by multiplying through both sides of system (1).

The choice of R is guided by the following definition:

$$R = (A_C^T A_C)^{-1} A_C^T \tag{3}$$

where,  $A^T$  denotes the transpose of A.

Assuming the system (1) is in preconditioned form,

$$Mx = r \tag{4}$$

where,  $M_{i,j} \in M$  and  $r_i \in [b]$ , we introduce the interval Cholesky method (see also (Alefeld and Mayer, 2007) in the form

$$[l]_{jj} = \left( [m]_{jj} - \sum_{k=1}^{j-1} [l]_{jk}^2 \right)^{1/2} \tag{5}$$

$$[l]_{ij} = \left( [m]_{ij} - \sum_{k=1}^{j-1} [l]_{ik} [l]_{jk} \right) / [l]_{jj} \tag{6}$$

$$i = j + 1, \dots, n, \quad j = 1, 2, \dots, n$$

$$[l]_i = \left( [y]_i - \sum_{j=i}^{i-1} [l]_{ij} \right) / [l]_{ii} \tag{7}$$

$$[x]_i^C = \left( [y]_i - \sum_{j=i+1}^n [l]_{ji} [x]_j^C \right) / [l]_{ii} \tag{8}$$

$$i = n, n-1, \dots, 1$$

It is assumed that  $0 < L_i$  ( $i=1,2,\dots,n$ ) and that interval Cholesky method is feasible.

**Definition 1:** (Alefeld and Mayer, 2007): Assuming that M has an interval Cholesky Factorization and given that, if either  $n = 1$ ,

$$[L] = \left( \sqrt{[m]_{11}} \right)$$

or  $n > 1$  and ,

$$[L] = \begin{pmatrix} \sqrt{[m]_{11}} & 0 \\ [c] / \sqrt{[m]_{11}} & [L]^1 \end{pmatrix}$$

we then define the interval Cholesky Schur complement by the equation:

$$\Sigma_{[m]}^C = [m]^1 - [c][c]^T / [m]_{11}$$

if  $n > 1$  and  $0 \notin [m]_{11}$

We call the pair  $([L] [L]^T)$  as the interval Cholesky factorization of  $\Sigma_{[m]}^C$ . Obtaining interval Cholesky factorization of M via interval Gaussian elimination algorithm IGA (M), it is proved by (Alefeld and Mayer, 2007) that:

$$\Sigma_{[m]}^C \subseteq \Sigma_{[m]}^G$$

whenever the interval matrix M:

$$\in [A] = \begin{pmatrix} [m]_{11} & [C]^T \\ [V] & [m]^1 \end{pmatrix}$$

with Schur complement:

$$\Sigma_{[m]}^G = [M]^1 - [V][C]^T / [m]_{11}$$

and  $[V][C]^T = [V][C]^T$  exists, where

$$\Sigma_{[m]}^C, \Sigma_{[m]}^G$$

signify Interval Cholesky Factorization and Interval Gaussian algorithm for interval matrix M.

The parallelism of interval Cholesky method emanates from parallel computation of disk inversion via the inclusion isotonicity of square root by using the procedure described in Petkovic (1989) and Gargantini (1980).

To this end, we define disk  $[M_{i,j}]$  as the parametric notation:

$$\left\{ [m]_{ij} - m_{ijc} < \eta_j \right\} \tag{9}$$

where  $m_{ijc}$  is the centre and  $\eta_j$  is the radius of the matrix  $M_{i,j}$ .

Assuming the disk does not contain the origin i.e.  $|m_{ijc}| > \eta_j$ , the square root of  $[m]_{i,j}$  is defined by the equation (Gargantini, 1976; 1980) given that  $R_1$  and  $R_2$  are the roots of the disk in the form:

$$R_i = \left\{ \sqrt{|m_{ijc}|} e^{i\theta/2}, \frac{\eta_j}{\sqrt{|m_{ijc}|} + \sqrt{||m_{ijc}| - \eta_j}} \right\} \tag{10}$$

$$R_j = \left\{ \sqrt{|m_{ijc}|} e^{i(\theta/2 + \pi)}, \frac{\eta_j}{\sqrt{|m_{ijc}|} + \sqrt{||m_{ijc}| - \eta_j}} \right\} \tag{11}$$

Where  $R_1$  and  $R_2$  are mutually disjoint and none of them contains the origin.

We can shorten notation for  $R_1$  and  $R_2$  for brevity in the form:

$$R_1 = \{ \psi_1, w_1 \}, R_2 = \{ -\psi_1, w_1 \}$$

The choice of the roots to be taken in the square root formulas expressed in Eq. (10) and (11) is guided by the procedures discussed in (Petkovic, 1989).

For the above square root Eq. in (10) and (11) to be effective, the centered disk inversion due to (Carstensen and Petkovic, 1994) has been introduced in the form:

$$m_{ij} = \{ \psi_1, w_1 \} = \left\{ \frac{1}{m_{ijc}}, \frac{2w_1}{|m_{ijc}|^2 - w_1^2} \right\} \quad (12)$$

$m_{ijc} > w_1$

Let us note that  $\theta$  appearing in Eq. (10) and (11) denotes the value of the argument  $m_{ij}$ .

The existence verification for  $\sum_{[m]}^C$  relies solely on the existence verification for IGA(M) whenever the matrix M is symmetric. This is also not surprising since the matrix M is an M-matrix.

Defining the matrix D by:

$$D = \text{diag}(L_{11}, L_{22}, \dots, L_{mm})$$

and since D has positive diagonal entries, we can decompose M in the form  $M = LL^T = (LD^{-1})(DL^T)$  where from again:

$$[M] \subseteq (LD^{-1})(DL^T)$$

**Comparison with Rohn's method:** The Rohn's method is implied by the following theorem:

**Theorem 2:** (Rohn, 1996)

Let R be an arbitrary nxm matrix, suppose  $x_0$  and  $d > 0$  be arbitrary n-vectors such that:

$$Gd + g < d \quad (13)$$

holds where,

$$G = |I - RA_c| + |R|\Delta \quad (14)$$

$$g = |R(A_c x_0 - b_c)| + |R|(\Delta |x_0| + \delta) \quad (15)$$

Then an inclusion for the linear interval system is given by:

$$x \subseteq [x_0 - d, x_0 + d] \quad (16)$$

The  $x_0$  is taken to be  $Rb_c$  where  $d_0$  may be assumed to be 0.

A variant of Rohn's method is the interval version proposed in (Rump, 1993). It is based on the work of (Krawczyk, 1969). Iteratively, Rohn's method (13) is written in the form:

$$d_{k+1} = Gd_k + g + f \quad (17)$$

where f is an arbitrary small vector. The vector f plays a very crucial role in the convergence of Rohn's theorem.

## RESULTS AND DISCUSSION

The numerical test is taken from (Popova, 2006) given as problem 1. Consider a parameterized system (1.1) of order 6x4 matrix

$$A = \begin{pmatrix} a & a+1 & a+2 & a \\ a & a+2 & a+3 & a+1 \\ a+1 & a+2 & a+3 & a+2 \\ a+2 & a+3 & a+4 & a+3 \\ a+3 & a+4 & a+5 & a+5 \\ a+4 & a+5 & a+6 & a+7 \end{pmatrix}$$

$$b = \begin{pmatrix} 3 \\ 1 \\ -2 \\ -3 \\ 0 \\ 0 \end{pmatrix}, \quad \epsilon = \begin{pmatrix} 1 \\ -1 \\ 1 \\ -1 \end{pmatrix}$$

We are requested to solve  $[A]x = b$ ,  $[A] = ([A(a)], |a \in [0.999, 1.001]])$

The following results were obtained by (Popova, 2006) using block Krawczyk's algorithm

$$[X] = ([0.8629, 1.1371], [-1.8273, -0.1727], [0.4924, 1.5076], [-1.2742, -0.7257])^T$$

When the same problem 1 above was solved with Parallel Rohn's method of Eq. (13), the following results were obtained:

$$[X] = ([0.999, 1.001], [-1.00997, -0.999003], [0.998997, 1.001003], [-1.001, -0.999003])^T$$

The Interval Cholesky Method represented in midpoint –radius form gave the following results:

Interval Cholesky Method

Midpoint X	Radius X
1	0
-1.013684744	0.028789846
1	0.009105466
-1	0.009113722

**CONCLUSION**

The interval widths produced by Rohn’s method of (13) are by far the narrowest than those obtained via interval Cholesky method. The block Krawczyk’s method by (Popova, 2006) used in the above example gave inclusion bounds whose widths have the largest values. It follows that Rohn’s algorithm is the best as far as these investigations are concerned.

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