# High-Performance Computing Fall 2022 Assignment 3

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## 0.1 Introduction

In this assignment, the objective is implementing the Manager Worker MPI model for computing maximum Eigenvalues of random matrices and finding the distribution of the Eigenvalues for a large number of matrices. For a symmetric matrix  $A[n \times n]$  with elements normally distributed. An Eigenvector v lies along the same line as Av:  $A\mathbf{v} = \lambda \mathbf{v}$ , the Eigenvalue is  $\lambda$ , The above equation can also be written as:  $(A - \lambda) = 0$  In a random matrix,  $\lambda$ max is the maximum Eigenvalue, and the distribution of such maximum Eigenvalues can be determined.

### What is Eigenvalue?

The eigenvalue is explained to be a scalar associated with a linear set of equations which, when multiplied by a nonzero vector, equals to the vector obtained by transformation operating on the vector.

The hardware that we run our tests on is:

CPU Information Model name:

CPUs is 8 - Intel(R) Core(TM)

Cache size is 6144 kb

CPU Max MHZ = 3900

CPU Min MHZ = 1200

Threads per core is 2 Core(s) per socket: 4

#### **RAM Information**

Memtotal is 16310212 kb Cached is 1797260 kb

## 0.2 Method

In this assignment, we investigate the speed up and efficiency of parallelization by computing the former distribution of  $\lambda$ max for large and small matrices using the Manager-Worker MPI model. In addition, we want to show whether the distribution of  $\lambda$ max is normal or not. In this study, I run the code on SHARCnet to show the distribution of large matrices.

## **0.2.1** Implementation

For the present study, we implemented a Manager-Worker system with the manager keeping track of the number of computations, receiving and collecting the outputs of each worker, and giving instructions to the ideal worker. In each Worker, random matrices are generated, Eigenvalues are computed, and results are sent back to the Master.

## 0.3 Psuedocode

#### **0.3.1** Worker Pseudocode

**Data:** Boolean (conv= converges), seed for PRNG or pseudorandom number generator, recvd\_tags: Receiving tag it means communication between worker and manager, Tag: sending tag, Status: fail-safe, Eig = Eigenvalues, A = Matix, ierr: For errors

MPI\_Bcast = give – integer and receive from manager to all worker (One to all)

Allcoate Matrix

**Loop** MPI\_reev = Receive the seed (for generating random number) from the manager (integer) and any worker can receive

recv\_tag status = checks that worker received seed

Makes matrix and calculates eigenvalues

If no convergence then "bad end"

If convergence then give tag = 0 otherwise tag = 1

MPI\_send = Send back maximum of Eigenvalues and also Eigenvalues should not be an integer

#### End loop

Deallocate matrix (because we are done with memory)

End subroutine worker

Subroutine making matrix (Use lapack, Lapack is Linear Algebra Package)

Call DGEEV (DGEEV computes for an N-by-N real nonsymmetric matrix A, the eigenvalues and, optionally, the left and/or right eigenvectors.)

Subroutine compute Eigenvalues (use lapack)

### 0.3.2 Manager Pseudocode

**Data:** num\_procs (number of threads), Seed (PRNG), ierr (for errors), recv = 0 (counter for Eigenvalues), Worker ( worker that we communicate with), exit tags, sent tag, ndat (number of Eigenvalues( $\lambda$ )), P (loop counter), failed,

**Initialize:** set matrix and number of eigenvalues

B\_cast - send matrix

Allocate eigs (number of eigenvalues)

Initializing tags

Write to file eigs

## Loop

If a worker is ready, make seed

If recvd, number of eigenvalues, n\_procs 2

Otherwise stop

Because every matrix is random and has its own seed

Give seed, check for done

Worker is busy

Once exit tags are sent

Finish

MPI\_Recv (receive buf (which is not a singular integer))

Can receive from anybody

Find out which worker sent out eigenvalues

Find out if convergent

If tag=0 then

recvd = recevd + 1

Add eigenvalues to list

Write data

Otherwise, increase the failed counter

Print results

Now make worker available

#### End loop

Deallocate list

Initialization data

n = size of the matrix

ndat = number of eigenvalues

End subroutine manager

#### 0.3.3 Main Pseudocode

Module global: because for MKL & MPI must be global variables

MKL\_VSL\_TYPE Routines for quasi-random number generation( Quasi-random sequences are functions from the positive integers to the unit hypercube.)

**Data**: n (number of matrix size must be an integer), exit tag, void = 0, proc\_num,num\_procs ( must be an integer) matrix splitting parameters,

End global

Get random seed, use the hard drive and iostat=istat(Must be an integer) (An IOSTAT value is a value assigned to the variable for the IOSTAT= specifier.), end

Get a random number and use MKL, called every time a matrix is filled (with open random stream)

If ierr is not equal to 0 then print returns flags

End if

Deallocate pseudo-random number stream

Pseudo-random number stream

**Main:** Wtime is the time and variable ok (everything is okay with code ?)

call MPI (make sure everything is set up property)

If proc\_num is equal to 0 then use MPI library clock (wtime)

If everything is good and proc\_num is equal to 0 call manage and worker.

Print walltime

call finalize

#### End main

Initialize MPI (check everybody is okay)

If ierr is not equal to 0 then

print MPI comm size failed

End IF

Mpi comm rank: returns the calling process's rank in the specified communicator. It's often necessary for a process to know its own rank.

IF ierr is not equal 0 then

Print mpi comm rank failed

Finalizes all code, Call MPI Barrier (It is Communicator), Call MPI Initialized (The functions MPI\_INIT and MPI\_FINALIZE are used to initiate and shut down an MPI computation,).

# 0.4 Experiment

In this experiment, we investigate the distribution of the maximum Eigenvalues at different matrix sizes and different Eigenvalue numbers. We compare wall time and then used PYTHON For plotting the data. The results are as follows:

Matrix size	Number of Eigenvalues	Walltime
100×100	600	0.47 sec
$400 \times 400$	15000	336 sec
$2000 \times 2000$	600	2052 sec
$2000 \times 2000$	10000	SHARCnet(Goerges data)
$5000 \times 5000$	1000	SHARCnet(Goerges data)
$6000 \times 6000$	1100	91413 sec

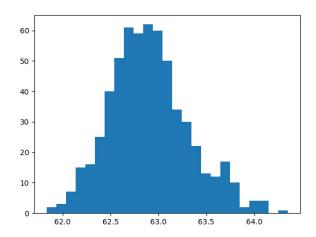


Figure 1:  $\lambda = 600$ , 1000x1000-sized matrices

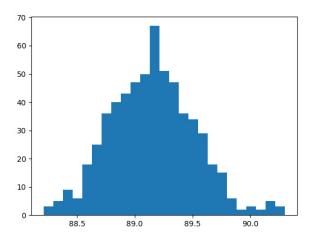


Figure 2:  $\lambda = 600$ , 2000x2000-sized matrices

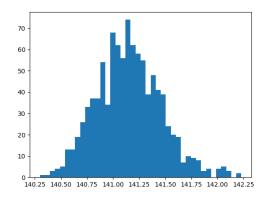


Figure 3:  $\lambda = 1000$ , 5000x5000-sized matrices

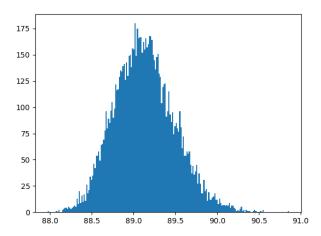


Figure 4:  $\lambda = 10000$ , 2000x2000-sized matrices

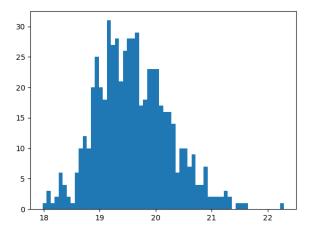


Figure 5:  $\lambda = 600$ , 100x100-sized matrices

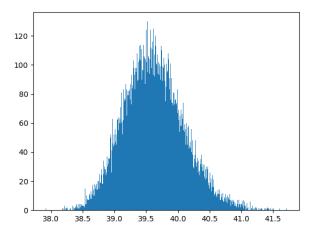


Figure 6:  $\lambda = 15000$ , 400x400-sized matrices

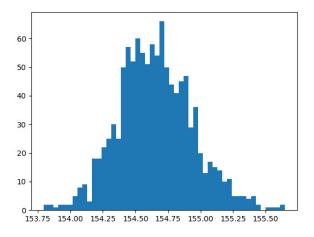
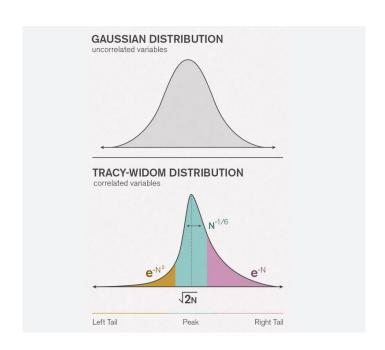


Figure 7:  $\lambda = 1100$ , 6000x6000-sized matrices



 $Figure~8:~\textit{Tracy-Widom~distribution~VS~gaussian~distribution}\\ https://www.wired.com/2014/10/tracy-widom-mysterious-statistical-law/$ 

In order to implement the code, Mersenne Twister pseudorandom number generation from the Intel Math Kernel Library (MKL) was used to seed the random numbers. LAPACK was implemented to calculate the maximum eigenvalues. The LAPACK subroutine is  $O(n^3)$ .

In large matrices, Lapack will use multi-threading automatically unless you tell it otherwise. Following that, we don't use multi-threading because it will affect on Wall times

generally, these results are presented in figures to show that the Eigenvalues follow a Tracy-Wisdom distribution with a Gaussian distribution as one limit for small sizes.

In these two figures 4 and 6 we can see that the distribution function which can be fitted to the matrices with small sizes would be similar to the Tracy-Widom distribution function presented in figure 8.

We can conclude that The fitted distribution function would be more similar to Tracy-Widom when the number of Max Eigenvalues increases.

#### The memory requirement is:

 $100 \times \text{N/MB}$ 

which N = Cpu size, M = matrix sizes - (for example  $1000 \times 1000$ ) B = 8 bytes.

This will show how much of the CPU is going to use. (percentage)

## 0.5 Conclusion

As a result of the computational experiments, we are able to conclude that the study's expected outcomes have been verified. When the number of Workers increased, MPI implementation of the Manager and Worker model resulted in a reduction in wallclock times. Well, it should model a Tracy-Widom distribution and not a Gaussian so we would reject the zero hypotheses.