

**MCS 572: Homework 2 (Revised)**  
**University of Illinois at Chicago (Professor Nicholls)**  
**Fall 2009**

Due Monday, October 12 by 2pm.

1. (a) Modify the matrix–vector multiplication code (`matvec.f`) to perform the Jacobi iteration for the solution of linear systems of equations  $A\vec{x} = \vec{b}$ .  
(b) Consider the matrix  $A \in \mathbf{R}^{n \times n}$  which is  $(n + 1)$  on the diagonal and 1 everywhere else. Pick the “exact solution”  $\vec{x}_{exact}$  which is all 1’s. What is the corresponding  $\vec{b}_{exact} = A\vec{x}_{exact}$  which gives this solution?  
For  $n = 10$  run the code (perform the Jacobi iteration with this  $(\vec{x}_{exact}, \vec{b}_{exact})$  pair until a relative error of  $10^{-10}$  is achieved) with  $N = 2, 4, 8$  processes. How many iterations are required? What speed–up do you notice?  
(c) Repeat part (b) with  $n = 100$ .  
(d) Repeat part (b) with  $n = 1000$ .
2. Exercises (Heath): 2.7, 2.21, 2.26, 2.31
3. Computer Problems (Heath): 2.2, 2.4, 2.6, 2.17
4. Exercises (Heath): 11.6, 11.7
5. Computer Problems (Heath): 11.12

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program main
include 'mpif.h'
integer MAX_ROWS, MAX_COLS, rows, cols
parameter (MAX_ROWS = 1000, MAX_COLS = 1000)
double precision a(MAX_ROWS,MAX_COLS), b(MAX_COLS), c(MAX_ROWS)
double precision buffer(MAX_COLS), ans

integer myid, master, numprocs, ierr, status(MPI_STATUS_SIZE)
integer i, j, numsent, sender
integer anstype, row

call MPI_INIT( ierr )
call MPI_COMM_RANK( MPI_COMM_WORLD, myid, ierr )
call MPI_COMM_SIZE( MPI_COMM_WORLD, numprocs, ierr )
master = 0
rows   = 100
cols   = 100

if ( myid .eq. master ) then
c  master initializes and then dispatches
c  initialize a and b (arbitrary)
do 20 j = 1,cols
    b(j) = 1
    do 10 i = 1,rows
        a(i,j) = i
10    continue
20    continue
    numsent = 0
c  send b to each slave process
    call MPI_BCAST(b, cols, MPI_DOUBLE_PRECISION, master,
&                MPI_COMM_WORLD, ierr)
c  send a row to each slave process; tag with row number
do 40 i = 1,min(numprocs-1,rows)
    do 30 j = 1,cols
        buffer(j) = a(i,j)
30    continue
    call MPI_SEND(buffer, cols, MPI_DOUBLE_PRECISION, i,
&                i, MPI_COMM_WORLD, ierr)
    numsent = numsent+1
40    continue
do 70 i = 1,rows
    call MPI_RECV(ans, 1, MPI_DOUBLE_PRECISION,
&                MPI_ANY_SOURCE, MPI_ANY_TAG,
&                MPI_COMM_WORLD, status, ierr)
    sender      = status(MPI_SOURCE)

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    anstype = status(MPI_TAG)      ! row is tag value
    c(anstype) = ans
    if (numsent .lt. rows) then    ! send another row
        do 50 j = 1,cols
            buffer(j) = a(numsent+1,j)
50          continue
            call MPI_SEND(buffer, cols, MPI_DOUBLE_PRECISION,
&                sender, numsent+1, MPI_COMM_WORLD, ierr)
            numsent = numsent+1
        else ! Tell sender that there is no more work
            call MPI_SEND(MPI_BOTTOM, 0, MPI_DOUBLE_PRECISION,
&                sender, 0, MPI_COMM_WORLD, ierr)
        endif
70    continue
    else
c        slaves receive b, then compute dot products until
c        done message received
        call MPI_BCAST(b, cols, MPI_DOUBLE_PRECISION, master,
&                MPI_COMM_WORLD, ierr)
c        skip if more processes than work
        if (rank .gt. rows)
&            goto 200
90    call MPI_RECV(buffer, cols, MPI_DOUBLE_PRECISION, master,
&                MPI_ANY_TAG, MPI_COMM_WORLD, status, ierr)
        if (status(MPI_TAG) .eq. 0) then
            go to 200
        else
            row = status(MPI_TAG)
            ans = 0.0
            do 100 i = 1,cols
                ans = ans+buffer(i)*b(i)
100          continue
            call MPI_SEND(ans, 1, MPI_DOUBLE_PRECISION, master,
&                row, MPI_COMM_WORLD, ierr)
            go to 90
        endif
200    continue
    endif

    call MPI_FINALIZE(ierr)
    stop
    end

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