

Homework Assignment 5 – due on Monday, November 6 (Midnight)

Description of Assignment:

Complete an MPI program(vecadd.c) that

- (i) decomposes A and B on p_0 to all other processors
- (ii) computes $C = A + B$ on all processors in parallel
- (iii) composes C on all other processors to p_0 .

<pre>#include <stdio.h> #include "mpi.h" #define N 24 main(int argc, char* argv[]) { int np, pid, i, tag = 0; float A[N], B[N], C[N]; MPI_Status status; MPI_Init(&argc, &argv); MPI_Comm_size(MPI_COMM_WORLD, &np); MPI_Comm_rank(MPI_COMM_WORLD, &pid); // initialization of A and B if (pid == 0) { for (i = 0; i < N; i++) { A[i] = i; B[i] = N-i; } } }</pre>	<pre>// (i) decomposition // (ii) addition for (i = 0; i < N/np; i++) C[i] = A[i] + B[i]; // (iii) composition // print results if (pid == 0) { for (i = 0; i < N; i++) printf("%2.1f ", C[i]); printf("\n "); } MPI_Finalize(); }</pre>
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How to proceed:

- (i) Use only MPI_Send() and MPI_Recv() for decomposition and composition.
- (ii) Run only 1, 2, 3, 4, 6, 12, 24 processors for tests.

Turnin the assignment:

After done your assignment, type **turnin** in your current working directory. You can retype the command at any time before the due date.