

# Homework 3

Due: Mar 15th @ 11:59 pm (as a typed, pdf file)

1. **OpenMP tasks:** Use OpenMP tasks to parallelize the following sequential code, you can assume that  $x$ ,  $y$ ,  $z$ ,  $f$ ,  $g$ ,  $h$ ,  $i$ , and  $j$  are all properly defined. Compute each of the variables  $a$ - $e$  using a task *specific* to that variable.

```
double x = ...;
double y = ...;
double z = ...;

double a = f(x);
double b = g(y);
double c = h(a,b);
double d = i(x,c);
double e = j(c,d,z);
```

We have provided starting code for you below:

```
double x = ...;
double y = ...;
double z = ...;

double a,b,c,d,e;

#pragma omp parallel
{

    // Compute a,b,c,d, and e in parallel using tasks.

}
```

2. **Pthreads:** Write a shared memory program using pthreads that computes the average of a distribution contained in array A in parallel by breaking the array A into smaller chunks. Each part a)-c) is marked in the comments for you to implement.

```
struct thread_info_struct {
    int start;    // start index of chunk for thread num
    int end;      // end index of chunk for thread num
    int* A;       // pointer to array
    pthread_t thread;
    int partial_sum;
}

void main() {
    /* assume thread attributes are initialized here */

    thread_info_struct thread_info[num_threads];
    int* A = (int*)malloc(1000*sizeof(int));
    assert (A != NULL);

    //initialize array
    for (int i = 0; i < 1000; i++) A[i] = rand() % 1000;

    for (int i = 0; i < num_threads; i++) {
        /* a) initialize thread_info_struct and create threads to compute
           partial_sum in parallel */

    }

    // join threads
    for (int i = 0; i < num_threads; i++) {
        pthreads_join(thread_info[i].thread, NULL);
    }

    double average = 0;
    /* b) compute average from partial sums */

    free(local_sum);
}

void* partial_sum(void* my_info) {
    /* c) Add your code here to compute local sum*/
}
```

3. **C++ Atomics:** Here is an incorrect implementation of a barrier, please explain where the error(s) is/are and what can go wrong if this is used.

```
std::atomic<int> x; // initialized to 0
int t = ...; // number of threads
void barrier() {
    int my_x = x.fetch_add(1);
    if (my_x == t) {
        x.store(0);
    } else {
        while (x.load() != t); // spin-wait
    }
}
```

(Hint: this implementation **will** work if only one barrier is needed throughout the entire program execution. Think about when multiple barriers are needed in the program.)

4. **MPI:** A tree has been constructed out of  $P$  processes such that process 0 is the root and the left and right children of a process  $i$  are processes  $2*i+1$  and  $2*i+2$  respectively. Each process contains an integer `my_val`.

Please use point-to-point communication **only** (i.e. send/rcv, no collectives) to implement the MPI reduction collective such that process 0's `sum_val` is the sum of each of the  $P$  process' `my_vals`. Your implementation should be equivalent to this MPI call:

```
MPI_Reduce(&my_val, &sum_val, 1, MPI_INT, MPI_SUM, 0,
MPI_COMM_WORLD);
```

```
// sum_val for process with rank 0 will be the sum of all my_val's
void my_reduce(int *my_val, int *sum_val) {
    int my_rank = ...; // assume already computed
    int num_ranks = ...; // assume already computed

    /* TODO: complete this function */
    int parent, leftchild, rightchild;

}
```