

DTS202TC Foundation of Parallel Computing

Lecture 4 OpenMP

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A1 Common Problems



FILE *f = fopen("test.txt", "r");

fscanf(f, "%d", &digit);
printf("%d \n", digit);

while ((count = fgetc(f)) != EOF) {

ungetc(count, f); // what does this do?

int digit, count:

fclose(f);

- Did not close the file after finishing the job (part of the coding quality)
- Did not use struct (part of the coding quality)
- One main function() contains everything
 - Very bad coding practice
- Hard-coding image array size
 int data[200000][200000];
- Output file name incorrect
- Makefile does not work, some students used cmake
- The parallel design should be in general, do not put it into one of the implementations.

Administrative



Week 5 CUDA

- 5 hours face-to-face lectures + lab from Nvidia Guest Lecturer.
- Same schedule as normal teaching, tutorial and lab for you to do exercise and test to get the CUDA certificate.
- Make sure you attend the lecture, you will be given a redeem code to redeem the course for free.
- · No recording due to the copyright.
- Individual Assessment 2 due on Saturday Dec. 23rd, 2023 @ 11:59pm
 - You must at least choose MPI.
- A2 FAQ
 - Can we use Gprof to get the elapsed time?
 - No, Gprof does not support multi-threading programs, please get the time manually by print it out in your code.
 - Can we make modifications of the serial code?
 - Yes, if you find problems of your A1 implementation, you can enhance it, and implement the parallel
 - Do we use the same pgm file for testing?
 - Yes, you can use the same pgm as A1 for your testing and evalution.

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Wrap up so far

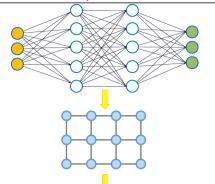


- Forster Method
 - Partitioning
 - Communication
 - Aggregation
 - Mapping
- Performance analysis
 - Speedup and Efficency
- Profiling
 - Gprof and other profiling tools
- Debugging C
 - Setting break points
 - Step over and step into

- Pthread
 - pthread_create, pthread_join
 - Critical Sections
 - Mutex
 - Semaphore
- MPI
 - Pointer-to-Point communication
 - MPI_Send and MPI_Receive
 - Collective communication
 - MPI_Reduce
 - MPI_Allreduce

Another Parallel Example

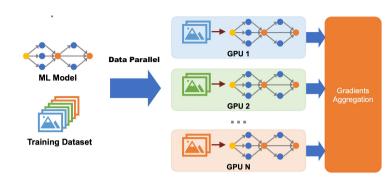




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Another Parallel Example – cont.



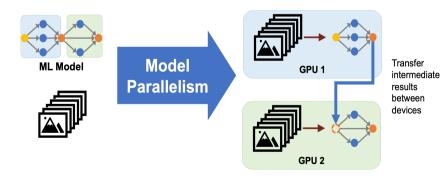


- Each GPU saves a replica of the entire model
- Cannot train large models that exceed GPU device memory

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Another Parallel Example – cont.





This week's topic



- OpenMP Basics
- Our First OpenMP Program
- Fundamental Concepts and Library Functions
- The Trapezoidal Rule
- Scope of Variables
- Task Parallelism
- The Reduction Clause
- The parallel for Directive

OpenMP



- An API for shared-memory parallel programming.
- MP = multiprocessing
- Designed for systems in which each thread or process can potentially have access to all available memory.

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Pragma



- Special preprocessor instructions.
- Typically added to a system to allow behaviors that aren't part of the basic C specification.
- Compilers that don't support the pragmas ignore them.

#pragma

OpenMP vs. Pthreads



- Pthreads requires that the programmer explicitly specify the behavior of each thread. OpenMP allows the compiler and run-time system to determine some of the details of thread behavior.
- Any Pthreads program can be used with any C compiler, provided the system has a Pthreads library. OpenMP requires compiler support for some operations, and hence it's entirely possible that you may run across a C compiler that can't compile OpenMP programs into parallel programs.
- Pthreads is lower level. Cost: Specify every detail of the behavior of each thread. OpenMP can be simpler to code some parallel behaviors. Cost: Some low-level thread interactions can be more difficult to program.

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"Hello, World" Using OpenMP



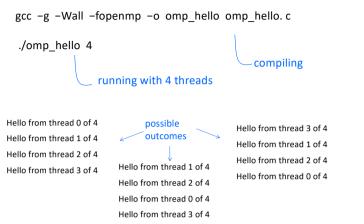
```
#include < stdio.h>
#include < stdlib.h>
#include <omp.h>
void Hello(void); /* Thread function */
int main(int argc, char* argv[]) {
  /* Get number of threads from command line */
  int thread count = strtol(argv[1], NULL, 10);
# pragma omp parallel num_threads(thread_count)
   Hello();
  return 0;
} /* main */
void Hello(void) {
  int my_rank = omp_get_thread_num();
  int thread_count = omp_get_num_threads();
  printf("Hello from thread %d of %d\n", my_rank, thread_count);
} /* Hello */
```

Compiling and Running



Demo





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Outline





- Our First OpenMP Program
- Fundamental Concepts and Library Functions
- The Trapezoidal Rule
- Scope of Variables
- Task Parallelism
- The Reduction Clause
- The parallel for Directive

OpenMP Syntax



- Most of the constructs in OpenMP are compiler directives
- #pragma omp directive [clause list]
- A parallel directive: # pragma omp parallel
- The number of threads that run the following structured block of code is determined by the run-time system.
- Function prototypes and types in the file
- #include <omp.h>



- There may be system-defined limitations on the number of threads that a program can start.
- The OpenMP standard doesn't guarantee that this will actually start thread count threads.
- Most current systems can start hundreds or even thousands of threads.
- Unless we're trying to start a lot of threads, we will almost always get the desired number of threads.

include <omp.h>

#ifdef _OPENMP

include <omp.h>

#endif

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In Case the Compiler doesn't Support OpenMP



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```
# ifdef _OPENMP
  int my_rank = omp_get_thread_num ( );
  int thread_count = omp_get_num_threads ( );
# else
  int my_rank = 0;
  int thread_count = 1;
# endif
```

OpenMP Library Functions



Control the number of threads and Processors:

```
#include <omp.h>
 void omp set num threads (int num threads);
 int omp get num threads ();
 int omp get max threads ();
 int omp get thread num ();
 int omp get num procs ();
 int omp_in_parallel ();
· Set and monitor thread creation:
 #include <omp.h>
 void omp set dynamic (int dynamic threads);
 int omp get dynamic ();
 void omp set nested (int nested);
 int omp get nested ();
Mutex:
 #include <omp.h>
 void omp_init_lock(omp_lock_t *lock);
 void omp destroy lock(omp lock t *lock);
 void omp set lock(omp lock t *lock);
 void omp unset lock(omp lock t *lock);
 int omp_test_lock(omp_lock t *lock);
• OpenMP also supports nested lock, which has similar semantics with simple lock.
```

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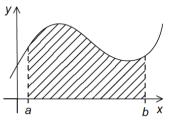
Outline

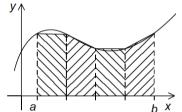


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The Trapezoidal Rule







If each subinterval has the same length h and if we define h=(b-a)/n, $x_i=a+ih$, $i=0,1,2,\ldots,n$, then our approximation will be

$$h[f(x_0)/2+f(x_1)+f(x_2)+\cdots+f(x_{n-1})+f(x_n)/2]$$

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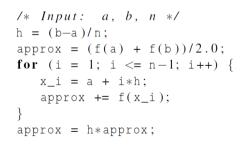
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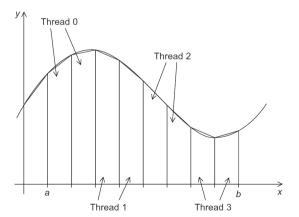
Serial Algorithm



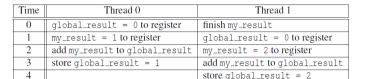
Assignment of Trapezoids to Threads











- Unpredictable results when two (or more) threads attempt to simultaneously execute: global result += my result;(critical section)
- Recall:

Race Condition, Critical Section



Critical Directive:

```
# pragma omp critical
 global result += my result;
                 only one thread can execute
```

the following structured block at a time

First OpenMP Trapezoidal Rule Program(1)



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```
#include < stdio . h>
#include < stdlib . h>
#include <omp.h>
void Trap(double a, double b, int n, double* global_result_p);
int main(int argc, char* argv[]) {
   double global_result = 0.0; /* Store result in global_result */
   double a, b;
                                /* Left and right endpoints
   int
                                /* Total number of trapezoids
         n;
          thread_count;
   thread_count = strtol(argv[1], NULL, 10);
   printf("Enter a, b, and n\n");
   scanf("%lf %lf %d", &a, &b, &n);
# pragma omp parallel num_threads(thread_count)
   Trap(a, b, n, &global_result);
   printf("With n = %d trapezoids, our estimate\n", n);
  printf("of the integral from %f to %f = %.14e\n",
     a, b, global_result);
   return 0:
} /* main */
```

First OpenMP Trapezoidal Rule Program(2)



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```
void Trap(double a, double b, int n, double* global_result_p) {
   double h, x, my_result;
   double local_a, local_b;
   int i, local_n;
   int my_rank = omp_get_thread_num();
   int thread_count = omp_get_num_threads();
   h = (b-a)/n;
   local n = n/thread count:
   local_a = a + my_rank*local_n*h;
   local_b = local_a + local_n*h;
   my_result = (f(local_a) + f(local_b))/2.0;
   for (i = 1; i \le local n - 1; i++)
    x = local_a + i*h;
     my_result += f(x);
   my_result = my_result*h;
# pragma omp critical
   *qlobal_result_p += my_result;
} /* Trap */
```

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- In serial programming, the scope of a variable consists of those parts of a program in which the variable can be used.
- In OpenMP, the scope of a variable refers to the set of threads that can access the variable in a parallel block.

Definition of Scope

Scope in OpenMP



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• A variable that can be accessed by all the threads in the team has shared scope.

- A variable that can only be accessed by a single thread has private scope.
- The default scope for variables declared before a parallel block is shared.

Example: Hello, World



```
#include < stdio.h>
#include < stdlib.h>
#include <omp.h>
void Hello(void); /* Thread function */
int main(int argc, char* argv[]) {
  /* Get number of threads from command line */
  int thread_count = strtol(argv[1], NULL, 10);
# pragma omp parallel num_threads(thread_count)
   Hello();
   return 0;
} /* main */
                     Private Scope
void Hello(void)
   int my_rank = omp_get_thread_num();
  int thread_count = omp_get_num_threads();
  printf("Hello from thread %d of %d\n", my_rank, thread_count);
} /* Hello */
```

Example: Trapezoidal Rule(1)



```
#include < stdio.h>
#include < stdlib . h>
#include <omp.h>
void Trap(double a, double b, int n, double* global_result_p);
int main(int argc, char* argv[]) {
   double | global result = 0.0; /* Store result in global_result */
   double a, b;
                                /* Left and right endpoints
                                /* Total number of trapezoids
          n:
          thread_count;
                            Shared Scope
   thread count = strtol(argv[1], NULL, 10);
   printf("Enter a, b, and n\n");
   scanf("%lf %lf %d", &a, &b, &n);
  pragma omp parallel num_threads(thread_count)
   Trap(a, b, n, &global_result);
   printf("With n = %d trapezoids, our estimate\n", n);
  printf("of the integral from %f to %f = %.14e\n",
     a, b, global_result);
   return 0;
} /* main */
```

Example: Trapezoidal Rule(2)



```
Private Scope
  void Trap(double a, double b, int n, double* global result p) {
    -double h, x, my result;
     double local_a, local_b;
     int i, local_n;
     int my_rank = omp_get_thread_num();
    int thread_count = omp_get_num_threads();
     h = (b-a)/n;
     local n = n/thread count:
     local_a = a + my_rank*local_n*h;
     local_b = local_a + local_n*h;
     my_result = (f(local_a) + f(local_b))/2.0;
     for (i = 1; i \le local_n - 1; i++) {
      x = local a + i*h;
       my result += f(x);
     my_result = my_result*h;
  # pragma omp critical
     *global_result_p += my_result;
 } _/* Trap */
                     Shared Scope
```

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Analysis of Trapezoidal Rule Program(1)



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In the OpenMP program, this more complex version is used to get global_result by adding each thread's local calculation.

```
void Trap(double a, double b, int n, double* global_result_p);
```

Although we'd prefer this for a serial implementation.

```
double Trap(double a, double b, int n);

qlobal result = Trap(a, b, n);
```



```
For the pointer version, we need to add each thread's local calculation to get global_result. If we use this, there's no critical section!

double Local trap(double a, double b, int n);
```

If we fix it like this, we force the threads to execute sequentially.

```
global_result = 0.0;
pragma omp parallel num_threads(thread_count)
{
    pragma omp critical
        global_result += Local_trap(double a, double b, int n);
}
```

We can avoid this problem by declaring a private variable inside the parallel block and moving the critical section after the function call.

```
global_result = 0.0;
# pragma omp parallel num_threads(thread_count)
{
    double my_result = 0.0; /* private */
    my_result += Local_trap(double a, double b, int n);
# pragma omp critical
    global_result += my_result;
}
```

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Syntax of Reduction Clause



- A reduction operator is a binary operation (such as addition or multiplication).
- A reduction is a computation that repeatedly applies the same reduction operator to a sequence of operands in order to get a single result.
- All of the intermediate results of the operation should be stored in the same variable: the reduction variable.

Reduction



A reduction clause can be added to a parallel directive.

```
global_result = 0.0;
# pragma omp parallel num_threads(thread_count) \
    reduction(+: global_result)

global_result += Local_trap(double a, double b, int n);
```

- When a variable is included in a reduction clause, the variable itself is shared. However, a private variable is created for each thread in the team.
- In the parallel block each time a thread executes a statement involving the variable, it uses the private variable. When the parallel block ends, the values in the private variables are combined into the shared variable.



• Start working on your individual assessment 2