Using OpenMP from C++

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Justification

OpenMP has the opportunity to exploit features of modern C++ that are not present in C. In this course we will explore:

- range-based iteration,
- differences in treatment between vectors and arrays, and various sophisticated reduction schemes.





Basic stuff





Bracket syntax

In keeping with the desire to get rid of the C Preprocessor (CPP) in C++, a new *syntax* for OpenMP *directives* was introduced:

```
1 // directive.cxx
2 int nthreads;
3 [[omp::directive( parallel ) ]]
4 [[omp::directive( master ) ]]
5 nthreads = omp_get_num_threads();
```





Output streams in parallel

The use of cout may give jumbled output: lines can break at each <<.
Use stringstream to form a single stream to output.





Parallel regions in lambdas

OpenMP parallel regions can be in functions, including lambda expressions.

```
1 const int s = [] () {
2   int s;
3  # pragma omp parallel
4  # pragma omp master
5   s = 2 * omp_get_num_threads();
6  return s; }();
```

('Immediately Invoked Function Expression')





Dynamic scope for class methods

Dynamic scope holds for class methods as for any other function:

```
Code:
1 // nested.cxx
2 class withnest {
3 public:
4 void f() {
   stringstream ss;
   << omp_get_num_threads()</pre>
   << '\n';
   cout << ss.str();</pre>
  };
11 };
12 int main() {
    withnest my_object;
14 #pragma omp parallel
15 my_object.f();
```

```
Output:

1 executing:

$\times OMP_MAX_ACTIVE_LEVELS=2$

$\times OMP_PROC_BIND=true$

$\times OMP_NUM_THREADS=2$

$\times ./nested$
2 2
3 2
```





Privatizing class members

Class members can only be privatized from (non-static) class methods.

In this example f can not be static:

```
1 // private.cxx
2 class foo {
3 private:
4   int x;
5 public:
6   void f() {
7  #pragma omp parallel private(x)
8   somefunction(x);
9  };
10 };
```

You can not privatize just a member:

```
1 // privateno.cxx
2 class foo { public: int x; };
3 int main() {
4    foo thing;
5  #pragma omp parallel private(thing.x) // NOPE
```





More privatization

Reference types can not be private.

Private objects are constructed/destructed by each thread.





Vectors are copied, unlike arrays, 1

C arrays: private pointer, but shared array:

```
Code:
1 // alloc.c
2 int *arrav =
    (int*) malloc(nthreads*sizeof(int));
4 for (int i=0: i<nthreads: i++)
   arrav[i] = 0;
7 #pragma omp parallel firstprivate(array)
8 {
    int t = omp_get_thread_num();
10 // ptr arith: needs private array
    array += t;
    arrav[0] = t:
12
13 }
14 // ... print the array
```

```
Output:

1 Array result:
2 0:0, 1:1, 2:2, 3:3,
```





Vectors are copied, unlike arrays, 2

C++ vectors: copy constructor also copies data:

```
Code:
1 vector<int> array(nthreads);
2 #pragma omp parallel firstprivate(array)
3 {
4   int t = omp_get_thread_num();
5   array[t] = t+1;
6 }
7 // ... print the array
```

```
Output:

1 Array result:
2 0:0, 1:0, 2:0, 3:0,
```





Parallel loops





Questions

- 1. Do regular OpenMP loops look different in C++?
- 2. Is there a relation between OpenMP parallel loops and iterators?
- 3. OpenMP parallel loops vs parallel execution policies on algorithms.





Range syntax

Parallel loops in C++ can use range-based syntax as of OpenMP-5.0:

```
1 // vecdata.cxx
2 vector<float> values(100);
3
4 #pragma omp parallel for
5 for ( auto& elt : values ) {
6   elt = 5.f;
7 }
8
9 float sum{0.f};
10 #pragma omp parallel for reduction(+:sum)
11 for ( auto elt : values ) {
12   sum += elt;
13 }
```

Tests show exactly the same speedup as the C code.





General idea

OpenMP can parallelize any loop over a C++ construct that has a 'random-access' iterator.





C++ ranges header

The C++20 ranges library is supported:





C++ ranges speedup

```
1 ==== Run range on 1 threads ===
2 sum of vector: 50000005000000 in 6.148
3 sum w/ drop 1: 50000004999999 in 6.017
4 sum times 2 : 100000010000000 in 6.012
5 ==== Run range on 25 threads ====
6 sum of vector: 50000005000000 in 0.494
7 sum w/ drop 1: 50000004999999 in 0.477
8 sum times 2 : 100000010000000 in 0.489
9 ==== Run range on 51 threads ====
10 sum of vector: 50000005000000 in 0.257
11 sum w/ drop 1: 50000004999999 in 0.248
12 sum times 2 : 100000010000000 in 0.245
13 ==== Run range on 76 threads ====
14 sum of vector: 50000005000000 in 0.182
15 sum w/ drop 1: 50000004999999 in 0.184
16 Sum times 2 : 100000010000000 in 0.185
17 ==== Run range on 102 threads ====
18 sum of vector: 50000005000000 in 0.143
19 sum w/ drop 1: 50000004999999 in 0.139
20 sum times 2 : 100000010000000 in 0.134
21 ==== Run range on 128 threads ====
22 sum of vector: 50000005000000 in 0.122
23 sum w/ drop 1: 50000004999999 in 0.11
24 sum times 2 : 100000010000000 in 0.106
```

25 scaling results in: range-scaling-ls6.out





Ranges and indices

Use iota_view to obtain indices:

```
1 // iota.cxx
2 vector<long> data(N);
3 # pragma omp parallel for
4 for ( auto i : std::ranges::iota_view( OUZ,data.size() ) )
5   data[i] = f(i);
```

Note that this uses C++23 suffix for unsigned size_t. For older versions:

```
1 iota_view( static_cast<size_t>(0),data.size() )
```





Custom iterators, 0

Recall that

Short hand:

for:

If we want

we need a sub-class for the iterator with methods such as begin, end, * and ++.

Probably also += and -





Custom iterators, 1

OpenMP can parallelize any range-based loop with a random-access iterator.

Class:

```
1 // iterator.cxx
2 template<typename T>
3 class NewVector {
4 protected:
5   T *storage;
6   int s;
7 public:
8 // iterator stuff
9   class iter;
10   iter begin();
11   iter end();
12 };
```

```
1 NewVector<float> v(s);
2 #pragma omp parallel for
3 for ( auto e : v )
4 cout << e << " ";</pre>
```

Main:





Custom iterators, 2

Required iterator methods:

```
1 NewVector<T>::iter& operator++();
2 T& operator*();
3 bool operator==( const NewVector::iter &other ) const;
4 bool operator!=( const NewVector::iter &other ) const;
5 // needed for OpenMP
6 int operator-
7 ( const NewVector::iter& other ) const;
8 NewVector<T>::iter& operator+=( int add );
```

This is a little short of a full random-access iterator; the difference depends on the OpenMP implementation.





Custom iterators, exercise

Write the missing iterator methods. Here's something to get you started.

```
1 template<typename T>
2 class NewVector<T>::iter {
3 private: T *searcher;
4 };
5 template<typename T>
6 NewVector<T>::iter::iter( T* searcher)
7 : searcher(searcher) {};
8 template<typename T>
9 NewVector<T>::iter NewVector<T>::begin() {
10 return NewVector<T>::iter(storage); };
11 template<typename T>
2 NewVector<T>::iter NewVector<T>::end() {
13 return NewVector<T>::iter(storage+NewVector<T>::s); };
```





Custom iterators, solution





Custom iterators, solution

```
1 template<typename T>
2 T& NewVector<T>::iter::operator*() {
3    return *searcher; };
4 // needed for OpenMP
5 template<typename T>
6 int NewVector<T>::iter::operator-
7    ( const NewVector<T>::iter& other ) const {
8    return searcher-other.searcher; };
```





OpenMP vs standard parallelism

Application: prime number marking (load unbalanced)

```
1 #pragma omp parallel for \
2    schedule(static)
3    for ( int i=0; i<nsize; i++) {
4        results[i] =
5        one_if_prime( number(i) );
6    }

1 // primepolicy.cpp
2 transform
3 ( std::execution::par,
4        numbers.begin(), numbers.end(),
5        results.begin(),
6    [] (int n ) -> int {
7        return one_if_prime(n); }
8    ):
```

Standard parallelism uses Threading Building Blocks (TBB) as backend





Timing

```
1 Threads: 1
2 TBR: Time:
                    392 msec
3 Stat: Time:
                    389 msec
4 Dyn: Time:
                    390 msec
 5
6 Threads: 25
 7 TRR: Time:
                      20 msec
                     32 msec
8 Stat: Time:
9 Dyn: Time:
                     17 msec
10
11 Threads: 51
12 TBB: Time:
                     13 msec
13 Stat: Time:
                     15 msec
14 Dyn: Time:
                      9 msec
15
16 Threads: 76
17 TBB: Time:
                      29 msec
18 Stat: Time:
                     11 msec
19 Dyn:
        Time:
                      6 msec
20
21 Threads: 102
22 TBB: Time:
                      61 msec
23 Stat: Time:
                      8 msec
24 Dyn: Time:
                      5 msec
25
26 Threads: 128
```

27 TBB:

Time:

80 msec 6 msec 4 msec



Reductions vs standard parallelism

Application: prime number counting (load unbalanced)

```
1 #pragma omp parallel for \
   schedule(guided,8) \
   reduction(+:prime_count)
4 for ( auto n : numbers ) {
   prime_count += one_if_prime( n );
6 }
1 // reducepolicy.cpp
2 prime count = transform reduce
    ( std::execution::par,
     numbers.begin(), numbers.end(),
     0,
     std::plus<>{},
     [] ( int n ) -> int {
         return one_if_prime(n); }
     ):
```





Timing

```
1 Threads: 1
2 TBR: Time:
                    391 msec
3 Stat: Time:
                    390 msec
4 Dyn: Time:
                    389 msec
 5
6 Threads: 25
 7 TRR: Time:
                      20 msec
8 Stat: Time:
                     17 msec
9 Dyn: Time:
                      17 msec
10
11 Threads: 51
12 TBB: Time:
                      13 msec
13 Stat: Time:
                      9 msec
14 Dyn: Time:
                      8 msec
15
16 Threads: 76
17 TBB: Time:
                      14 msec
18 Stat: Time:
                      8 msec
19 Dyn:
        Time:
                      5 msec
20
21 Threads: 102
22 TBB: Time:
                     76 msec
23 Stat: Time:
                      5 msec
24 Dyn: Time:
                      4 msec
25
26 Threads: 128
```

27 TBB:

Time:

80 msec 4 msec 3 msec



Reductions





Questions

- 1. Are simple reductions the same as in C?
- 2. Can you reduce std::vector like an array?
- 3. Precisely what can you reduce?
- 4. Any interesting examples?
- 5. Compare reductions to native C++ mechanisms.





Scalar reductions

Same as in C, you can now use range syntax for the loop.





Reductions on vectors

Use the data method to extract the array on which to reduce. However, this does not work:

```
1 vector<float> x;
2 #pragma omp parallel reduction(+:x.data())
```

because the reduction clause wants a variable, not an expression, for the array, so you need an extra bare pointer:

```
1 // reductarray.cxx
2 vector<int> data(nthreads,0);
3 int *datadata = data.data();
4 #pragma omp parallel for schedule(static,1) \
5 reduction(+:datadata[:nthreads])
```





Reduction on class objects

Reduction can be applied to any class for which the reduction operator is defined as <code>operator+</code> or whichever operator the case may be.

```
1 // reductclass.cxx
                                           1 vector< Thing >
2 class Thing {
                                               things(500, Thing(1.f));
3 private:
                                           3 Thing result(0.f);
4 float x{0.f}:
                                           4 #pragma omp parallel for \
                                                 reduction( +: result )
5 public:
    Thing() = default;
                                           6 for (const auto& t: things)
   Thing(float x): x(x) {};
                                           7 result = result + t:
   Thing operator+
        ( const Thing& other ) {
    return Thing( x + other.x );
11 }:
12 };
```

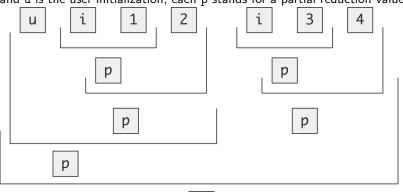
A default constructor is required for the internally used init value; see figure 34.





Reduction illustrated

Reduction of four items on two threads. i is the OpenMP initialization, and u is the user initialization; each p stands for a partial reduction value.









User-defined reductions, syntax

```
    #pragma omp declare reduction
    ( identifier : typelist : combiner )
    [initializer(initializer-expression)]
```





Reduction over iterators

Support for *C++ iterators*





Lambda expressions in declared reductions

You can use lambda expressions in the explicit expression for a declared reduction:

```
1 // reductexpr.cxx
2 #pragma omp declare reduction\
3    (minabs : int : \
4     omp_out = \
5          [] (int x,int y) -> int { \
6          return abs(x) > abs(y) ? abs(y) : abs(x); } \
7          (omp_in,omp_out) \
8     initializer (omp_priv=limit::max())
```

You can not assign the lambda expression to a variable and use that, because <code>omp_in/omp_out</code> are the only variables allowed in the explicit expression.





Example category: histograms

Count which elements fall into what bin:

```
for ( auto e : some_range )
histogram[ value(e)]++;
```

Collisions are possible, but unlikely, so critical section is very inefficient





Histogram: intended main program

Declare a reduction on a histogram object; each thread gets a local map:

```
1 /*
2 * Reduction loop in main program
3 */
4 bincounter<char> charcount;
5 #pragma omp parallel for reduction(+ : charcount)
6 for ( int i=0; i<text.size(); i++ )
7 charcount.inc( text[i] ):</pre>
```

Q: why does the inc not have to be atomic?





Histogram solution: reduction operator

Give the class a += operator to do the combining:

```
1 // charcount.cxx
2 template<typename key>
3 class bincounter : public map<key,int> {
4 public:
5 // merge this with other map
    void operator+=
        ( const bincounter<key>& other ) {
      for ( auto [k,v] : other )
        if ( map<key,int>::contains(k) )
          this->at(k) += v:
10
11
        else
          this->insert( {k,v} );
12
13
14 // insert one char in this map
    void inc(char k) {
      if ( map<kev.int>::contains(k) )
        this->at(k) += 1:
17
      else
19
        this->insert(\{k,1\}):
    };
20
21 }:
```





Histogram in native C++

Use atomics because there is no reduction mechanism:





Histogram in native C++, comparison

OpenMP reduction on array<int,26>:

```
1 Using atomics on 1 threads: time= 20.19 msec
2 OpenMP reduction on 1 threads: time= 1.966 msec
3 Using atomics on 5 threads: time= 315.855 msec
4 OpenMP reduction on 5 threads: time= 0.52 msec
5 Using atomics on 10 threads: time= 91.968 msec
6 OpenMP reduction on 10 threads: time= 0.364 msec
7 Using atomics on 30 threads: time= 249.171 msec
8 OpenMP reduction on 30 threads: time= 0.556 msec
9 Using atomics on 50 threads: time= 164.177 msec
10 OpenMP reduction on 50 threads: time= 0.904 msec
```





Exercise: mapreduce

Make an OpenMP parallel version of:

```
1 intcounter primecounter;
2 for ( auto n : numbers )
3    if ( isprime(n) )
4         primecounter.add(n);
```

where primecounter contains a map<int,int>.

Use skeleton: mapreduce.cxx





Example category: list filtering

The sequential code is as follows:

```
1 vector<int> data(100);
2 // fil the data
3 vector<int> filtered;
4 for ( auto e : data ) {
5    if ( f(e) )
6    filtered.push_back(e);
7 }
```





List filtering, solution 1

Let each thread have a local array, and then to concatenate these:

```
1 #pragma omp parallel
2 {
3   vector<int> local;
4 # pragma omp for
5   for ( auto e : data )
6   if ( f(e) ) local.push_back(e);
7   filtered += local;
8 }
```

where we have used an append operation on vectors:

```
1 // filterreduct.cxx
2 template<typename T>
3 vector<T>& operator+=( vector<T>& me, const vector<T>& other ) {
4  me.insert( me.end(), other.begin(), other.end() );
5  return me;
6 };
```





List filtering, not quite solution 2

We could use the plus-is operation to declare a reduction:

```
1 #pragma omp declare reduction\
2 ( \
3          +:vector<int>:omp_out += omp_in \
4          ) \
5     initializer( omp_priv = vector<int>{} )
```

Problem: OpenMP reductions can not be declared non-commutative, so the contributions from the threads may not appear in order.

```
Code:

1 #pragma omp parallel \
2 reduction(+: filtered)
3 {
4 vector<int> local;
5 # pragma omp for
6 for ( auto e : data )
7 if ( f(e) )
8 local.push_back(e);
9 filtered += local;
10 }
```

```
Output:

1 Mod 5: 80 85 90 95 100

→5 10 15 20 25 30

→35 40 45 50 55

→60 65 70 75
```





List filtering, task-based solution

Parallel region, without for:

```
Code:

1 // filtertask.cxx
2 vector<int> filtered;
3 int ithread=0;
4 #pragma omp parallel
5 {
6 vector<int> local;
7 int threadnum = omp_get_thread_num();
8 # pragma omp for
9 for ( auto e : data )
10 if ( e%5==0 )
11 local.push_back(e);
12 // create task to add local to filtered
```





List filtering, task-based solution'

The task spins until it's its turn:

```
Code:

1 # pragma omp task \
2     shared(filtered,ithread)
3     {
4 // wait your turn
5     while (threadnum>ithread) {
6 # pragma omp taskyield
7     }
8 // merge
9     filtered += local;
10     ithread++;
11 }
```





Templated reductions

You can reduce with a templated function if you put both the declaration and the reduction in the same templated function:

```
1 template<typename T>
2 T generic_reduction( vector<T> tdata ) {
3 #pragma omp declare reduction
4 (rwzt:T:omp_out=reduce_without_zero<T>(omp_out,omp_in))
5 initializer(omp_priv=-1.f)
6
7 T tmin = -1;
8 #pragma omp parallel for reduction(rwzt:tmin)
9 for (int id=0; id<tdata.size(); id++)
10 tmin = reduce_without_zero<T>(tmin,tdata[id]);
11 return tmin;
12 };
```

which is then called with specific data:

```
1 auto tmin = generic_reduction<float>(fdata);
```





More topics





Threadprivate random number generators

The new C++ random header has a threadsafe generator, by virtue of the statement in the standard that no STL object can rely on global state. The usual idiom can not be made threadsafe because of the initialization:

```
1 static random_device rd;
2 static mt19937 rng(rd);
```

However, the following works:

```
1 // privaterandom.cxx
2 static random_device rd;
3 static mt19937 rng;
4 #pragma omp threadprivate(rd)
5 #pragma omp threadprivate(rng)
6
7 int main() {
8
9 #pragma omp parallel
10 rng = mt19937(rd());
```





Threadprivate random use

Based on the previous note, you can use the generator safely and independently:

```
1 #pragma omp parallel
2  {
3    stringstream res;
4    uniform_int_distribution<int> percent(1, 100);
5    res << "Thread " << omp_get_thread_num() << ": " << percent(rng) << "\n";
6    cout << res.str();
7  }</pre>
```





Uninitialized containers

Multi-socket systems:

parallel initialization instantiates pages on sockets: 'first touch'

```
double *x = (double*)malloc( N*sizeof(double));
#pragma omp parallel for
for (int i=0; i<N; i++)
x[i] = f(i);</pre>
```

This does not work with

```
std::vector<double> x(N);
#pragma omp parallel for
for (int i=0; i<N; i++)
x[i] = f(i);</pre>
```

because of value initialization in the vector container.





Uninitialized containers, 2

Trick to create a vector of uninitialized data:

```
1 // heatalloc.cxx
2 template<typename T>
3 struct uninitialized {
4  uninitialized() {};
5   T val;
6   constexpr operator T() const {return val;};
7   T operator=( const T&& v ) { val = v; return val; };
8 };
```

so that we can create vectors that behave normally:

```
1 vector<uninitialized<double>> x(N),y(N);
2
3 #pragma omp parallel for
4 for (int i=0; i<N; i++)
5  y[i] = x[i] = 0.;
6 x[0] = 0; x[N-1] = 1.;</pre>
```

(Question: why not use reserve?)





Uninitialized containers, 3

Easy way of dealing with this:

```
1 template<typename T>
2 class ompvector : public vector<uninitialized<T>> {
3 public:
4    ompvector( size_t s )
5    : vector<uninitialized<T>>::vector<uninitialized<T>>(s) {};
6 };
```





Atomic scalar updates

Can you atomically update scalars?

- Make an object that has data plus a lock;
- Disable copy and copy-assignment operators;
- Destructor does omp_destroy_lock;
- Overload arithmetic operator.

(Quick self-test: why lock, not critical?)





Atomic updates: class with OMP lock

```
1 // lockobject.cxx
2 class atomic_int {
3 private:
    omp_lock_t the_lock;
    int _value{0};
6 public:
    atomic int() {
      omp_init_lock(&the_lock);
    atomic_int( const atomic_int& )
10
        = delete;
11
    atomic_int& operator=( const atomic_int& )
12
        = delete;
13
    ~atomic int() {
14
      omp_destroy_lock(&the_lock);
15
    };
16
```





Atomic updates: atomic ops

```
1 int operator +=( int i ) {
2 // atomic increment
3 omp_set_lock(&the_lock);
4 _value += i; int rv = _value;
5 omp_unset_lock(&the_lock);
6 return rv;
7 };
```





Atomic updates: usage

```
1 atomic_int my_object;
2 vector<std::thread> threads;
3 for (int ithread=0;
4    ithread<NTHREADS;
5    ithread++) {
6    threads.push_back
7    ( std::thread(
8      [=,&my_object] () {
9         for (int iop=0; iop<nops; iop++)
10         my_object += 1; } ) );
11 }
12 for ( auto &t : threads )
13    t.join();</pre>
```





Atomic updates, comparison to native

Timing comparison on simplest case:

```
Object with built-in lock:
                                              1 std::atomic<int> my_object{0};
                                              2 #pragma omp parallel for
 1 atomic_int mv_object;
                                              3 for ( size_t update=0;
2 vector<std::thread> threads:
                                                      update<NTHREADS*nops:
 3 for (int ithread=0;
                                                      update++) {
       ithread<NTHREADS:
                                                 mv_object += 1;
       ithread++) {
   threads.push_back
                                              8 result = mv_object;
      ( std::thread(
        [=,&mv_object] () {
           for (int iop=0; iop<nops; iop++)</pre>
             my_object += 1; } ) );
10
11 }
12 for ( auto &t : threads )
    t.join();
```

Native C++ atomics:

Native solution is 10x faster.





False sharing prevention

```
#include <new>
1
      #ifdef __cpp_lib_hardware_interference_size
      const int spread = std::hardware_destructive_interference_size
              / sizeof(datatype);
      #else
      const int spread = 8;
      #endif
8
      vector<datatype> k(nthreads*spread);
10
      #pragma omp parallel for schedule( static, 1 )
11
      for ( datatype i = 0; i < N; i++ ) {
12
        k[ (i%nthreads) * spread ] += 2;
13
14
```

Since C++17





Beware vector-of-bool!

Does not compile:

```
1 // boolrange.cxx
2 vector<bool> bits(1000000);
3 for ( auto& b : bits )
4  b = true;
```

More subtle:





CMake



