# Into to C/C++ pragma-based parallelism

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# We will mostly focus on C

- "Full" C++ hides a lot of details from the final user
  - In C, you explicitly tell how to do things, step by step
  - C better for understanding how things work
- Most of the presented concepts still apply to "full" C++
  - But you may need to understand the minute details that the compiler normally tries to hide from you!
  - Wherever appropriate, will call out known "gotchas"

### We will focus on portable methods

- Life is too short to re-implement your apps for every architecture!
  - That's why we use a compiler and not ASM to start with!
- Picking a vendor specific toolkit may give you better performance
   But with dominant vendors changing, hard to maintain/port
  - CPUs pthreads
  - NVIDIA GPUs CUDA
  - AMD GPUs ROCm/HIP
  - Intel GPUs OneAPI/SYCL

# We will focus on portable methods

• Life is too short to re-

No magic solutions, some porting/tuning always needed with new platforms.

- NVIDIA GPUS
- AMD GPUs ROCm/HIP
- Intel GPUs OneAPI/SYCL

capps for every architecture! start with!

> better performance to maintain/port

# We will focus on portable methods

- Life is too short to re-implement your apps for every architecture!
  - That's why we use a compiler and ASM to start with!
- Picking a vendor specification
   But with dominant
  - CPUs pthreads
  - NVIDIA GPUs CU
  - AMD GPUs ROCm/HIP
  - Intel GPUs OneAPI/SYCL

Will focus on mature solutions, so you can start using them now.

(Some cutting-edge solutions do promise to make your life easier)

### Virtually all parallelism in loops

- Compilers have hard time parallelizing sequential code
  - Not completely impossible, but very limited potential
  - Sometimes you can help
    - But often makes code hard to maintain

```
a1 += 2*b1;
a2 += 2*b2;
a3 += 2*b3;
a4 += 2*b4;
```

- Loops are the natural concept for expressing parallelism, if
  - Each iteration independent
  - Number of iterations known in advance

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- Loops are the natural concept for expressing parallelism, if
  - Each iteration independent
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```
for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
}
```

A good compiler will issue vectorized code for both snippets

### Loops and dependencies

• Loops are the natural concept for expressing parallelism, if

• Each iteration independent

Number of iterations known in advance

```
for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
}
```

As long as A and B do not overlap, it is pretty obvious each iteration is independent.

### Dependencies hard to infer

- Loops are the natural concept for expressing parallelism, if
  - Each iteration independent
  - Number of iterations known in advance

- Doesn't take much for compiler to give up
  - 1st priority for the compiler is to generate correct code
  - Optimization/speed is distant 2nd

```
for (int i=0; i<N; i++) {
    A[i] += 3*myfunc(A,B,i)
}
```

Good luck, if **myfunc** not inlined and/or not trivial.

### Dependencies hard to infer

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Not to mention "smart" users

```
for (int i=0; i<N; i++) {
    A[i] += 3*myfunc(A,B,i)
}
```

```
for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
    if(A[i]>3) N=A[i];
}
```

# Helping the compiler with pragmas

- OpenMP has emerged as the major pragma-based parallelization paradigm
  - Especially on CPUs
  - Just decorate the loop with a pragma
- Explicitly tell the compiler that parallelization is both possible and desirable

```
#pragma omp parallel for
for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
}</pre>
```



https://www.openmp.org

```
#pragma omp parallel for
for (int i=0; i<N; i++) {
    A[i] += 3*myfunc(A,B,i)
}</pre>
```

# Helping the compiler with pragmas

 OpenMP parallel Compiler will trust you! ttps://www.openmp.org If the code was not independent, you will get unexpected results. • Explicity is both possible and desire #pragma omp parallel for **#pragma omp parallel for** for (int i=0; i<N; i++) { for (int i=0; i<N; i++) { A[i] += 3\*B[i]A[i] += 3\*myfunc(A,B,i)

#### **GPU** extensions

- OpenMP recently added GPU support
  - The notion of "OMP TARGET"
- Slightly different syntax, but similar in concept

```
#pragma omp target teams distribute parallel for simd
for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
}</pre>
```



https://www.openmp.org

### **GPU pragmas - OpenACC**

OpenACC was the pioneer in GPU-based pragma-based parallelization



#pragma acc parallel loop gang vector
for (int i=0; i<N; i++) {
 A[i] += 3\*B[i]
}</pre>



https://www.openacc.org



Much easier to use than OpenMP Target in my opinion.

#### Remember the launch cost

- Every parallel section incurs a "launch cost"
  - Very expensive on GPUs
  - But not negligible on CPUs, either
- Pack as much as possible in a single parallel section
  - Can be order of magnitude faster

```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     A[i] += 3*B[i]
  }
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     C[i] += 4*D[i]
  }</pre>
```



```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
    C[i] += 4*D[i]
}</pre>
```

#### But don't break semantics!

- Every parallel section incurs a "launch cost"
  - Very expensive on GPUs
  - But not negligible on CPUs, either
- Pack as much as possible in a single parallel section
  - Can be order of magnitude faster
  - But make sure you preserve the independence of the iterations!

```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     A[i] += 3*B[i]
  }
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     C[i] += 4*A[1+(2*i+5)%N]
  }</pre>
```



```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
    A[i] += 3*B[i]
    C[i] += 4*A[1+(2*i+5)%N]
}</pre>
```

## Remember memory locality

- Do not iterate over same buffer multiple times
  - If at all possible (dependencies)
- Pack as much as possible in a single parallel section
  - Also use temp variables when appropriate
  - Will be order of magnitude faster on large buffers

```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     A[i] += 3*B[i]
  }
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
     C[i] += 4*A[i] + 3*B(i)
  }</pre>
```

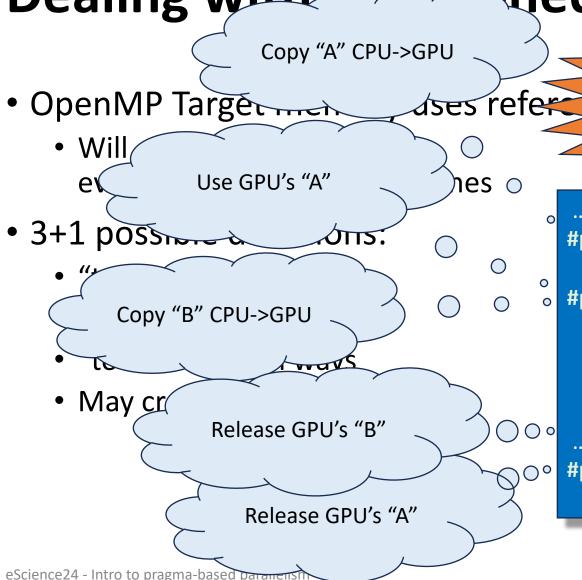


```
#pragma omp target teams \
  distribute parallel for simd
  for (int i=0; i<N; i++) {
    b3 = 3*B[i]
    a2 = A[i] + b3
    A[i] = a2
    C[i] += 4*a2 + b3
}</pre>
```

- Remainder: CPU and GPU memories are (typically) independent
- By default, OpenMP will copy all buffers
  - From CPU to GPU memory before GPU kernel start
  - From GPU to CPU memory after GPU kernel completion
- If buffers only needed on one side, you can avoid the transfer
  - Resulting in drastic speedup!

- OpenMP Target memory uses reference counters
  - Will only transfer once,
     even if mentioned several times
- 3+1 possible directions:
  - "to" CPU ->GPU
  - "from" GPU -> CPU
  - "tofrom" both ways
  - May create uninitialized buffer

Copy "A" CPU->GPU • OpenMP Target .... dses reference counters Wil Use GPU's "A" nes o • 3+1 possione 0 Copy "B" CPU->GPU May cr Release GPU's "B"  $\bigcirc \circ$ Release GPU's "A" eScience24 - Intro to pragma-based para



In this example, CPU's "A" never updated

# Ready for some hands-on?