Lecture #16 - Adaptive Quadrature and Load Balancing

AMath 483/583

Announcements

- Homework #4 Due Next Friday @ 5:00pm (Last Day of Class)
 - short, but conceptually challenging
 - (i.e. not something you can do at the last minute)
 - clarification in this lecture
- Today:
 - MPI on multiple machines
 - Return to OpenMP "nested parallelism"

Quick Aside

MPI is for distributed memory environments

```
$ mpiexec -n 4 -H [server list] ./myprog
```

Can also create a hostfile:

```
# my_hostfile
foo.example.com
bar.example.com slots=2
foobar.example.com slots=4 max-slots=4
```

Quick Aside

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# my_hostfile
foo.example.com
bar.example.com slots=2
foobar.example.com slots=4 max-slots=4
```

Quad-core machine - ensure no more than 4 processes run here

Disclaimer

- Hardest part is setting up computers / ssh-keys
- In following demo
 - americano.amath.washington.edu can send requests to mocha.amath.washington.edu
 - not vice versa (?)

Demo

hello-distributed.c

Clarify some misconceptions: use the <u>strategy</u>
 from shift.c, not the same algorithm

```
double* u;
```

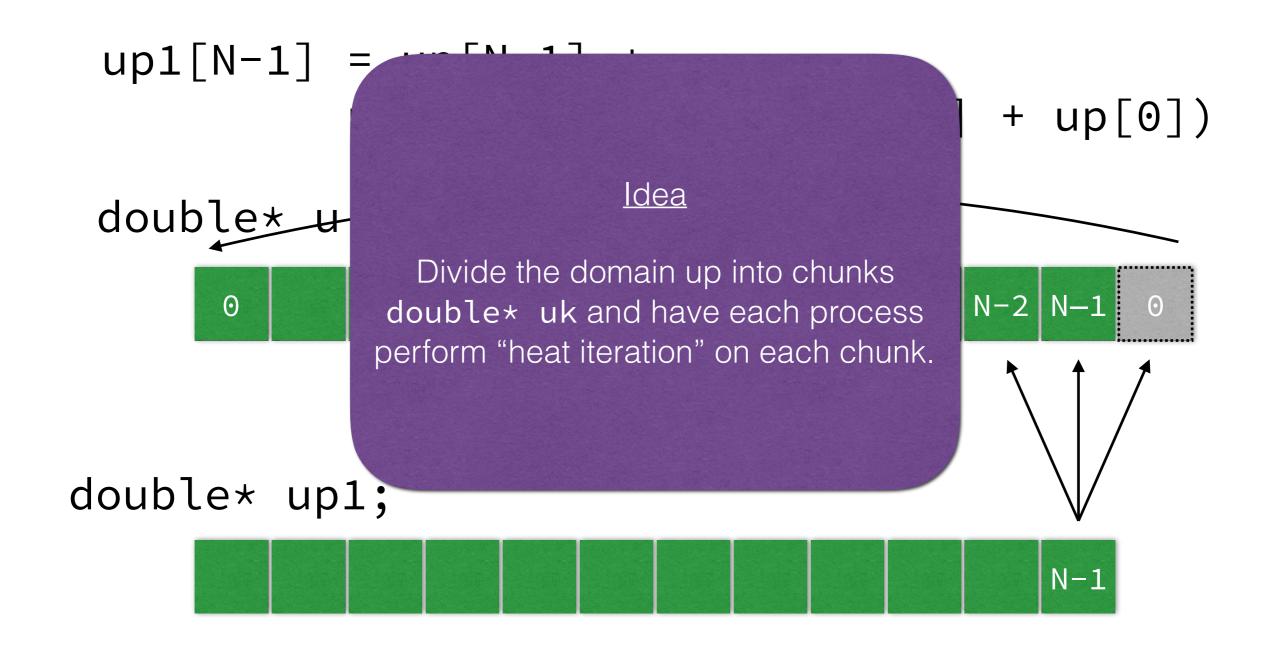


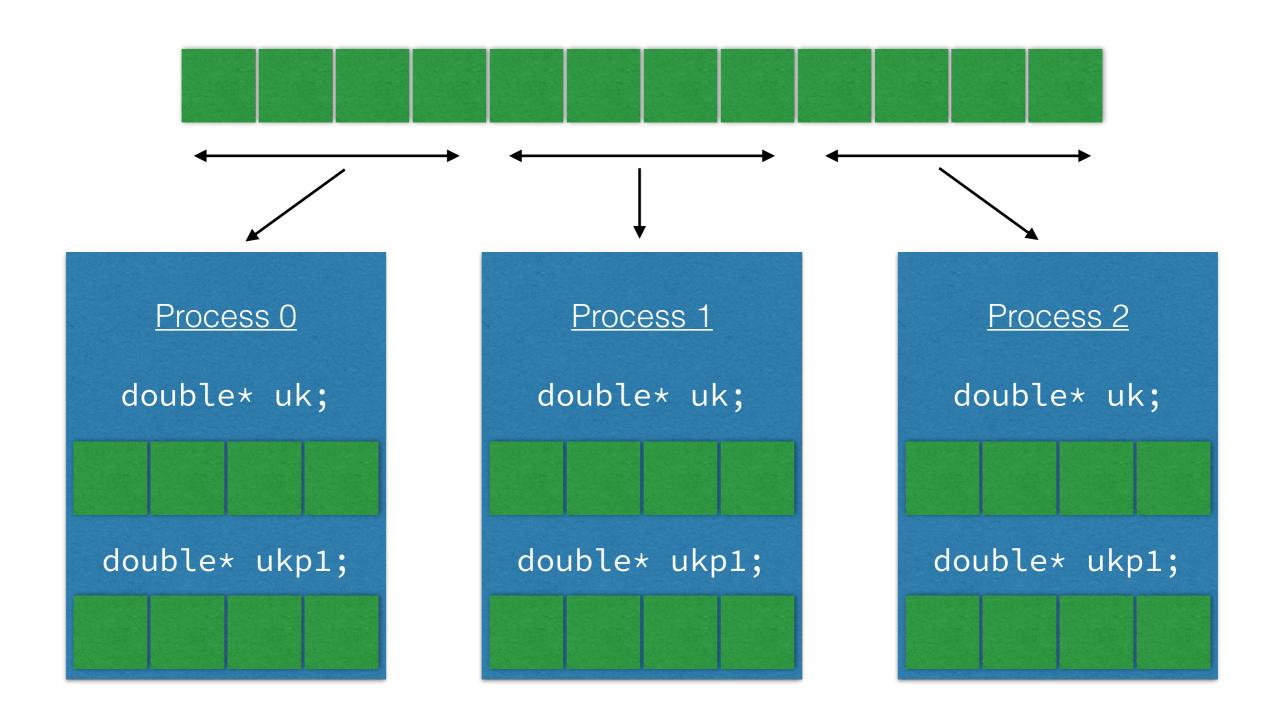
```
double* up1;
```

```
up1[i] = up[i] +
          nu*(up[i-1] - 2*up[i] + up[i+1])
 double* u;
double* up1;
```

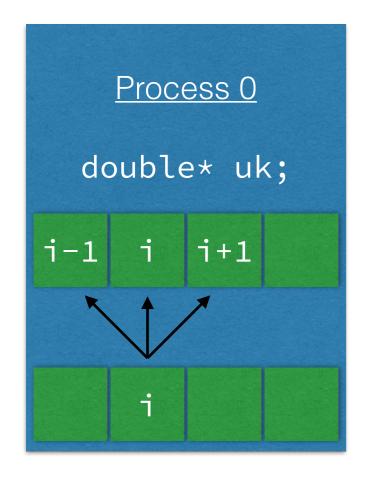
Solving the periodic

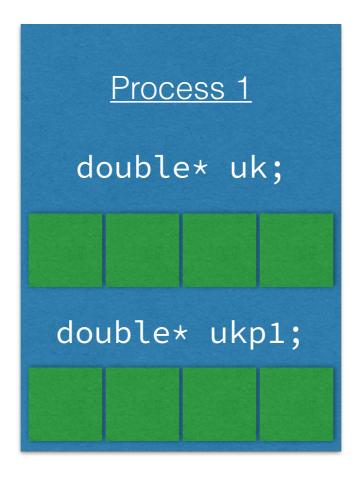
```
domain problem.
 up1[N-1] = up[N-1] +
           nu*(up[N-2] - 2*up[N-1] + up[0])
 double*
                                      N-2 N-1
      0
double* up1;
```

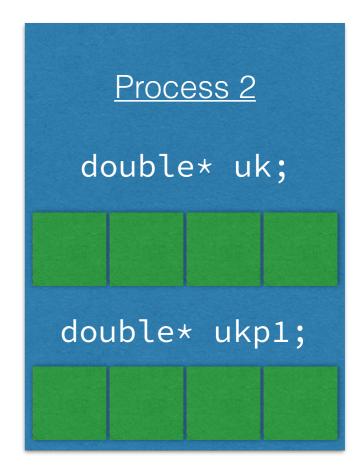




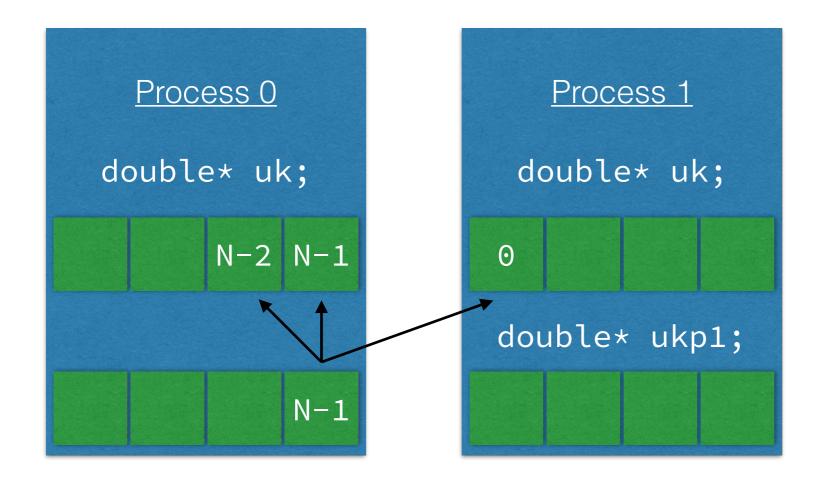
```
ukp1[i] = ukp[i] +
nu*(ukp[i-1] - 2*ukp[i] + ukp[i+1])
```

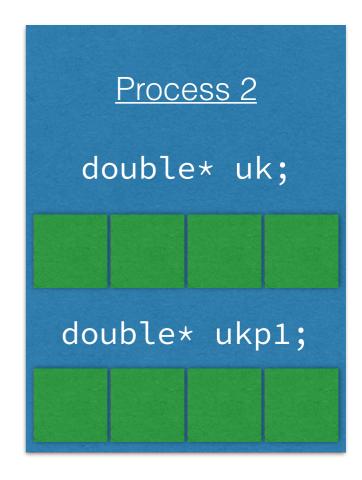






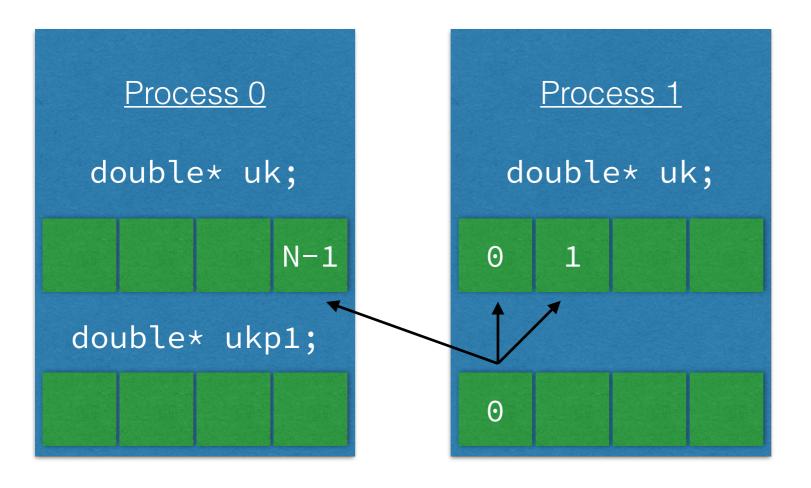
• **Uh oh**. Need to get uk information from neighboring processes. (Send / Recv)

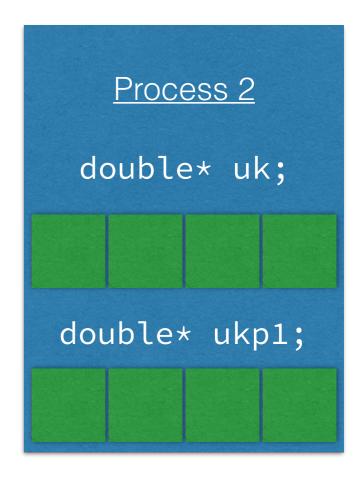




• **Same Issue**. Left side of process need info from left proc. Right side of process needs info from right proc.

(Hint: multiple ghost cells. More than one. More than two?)





Adaptive Quadrature

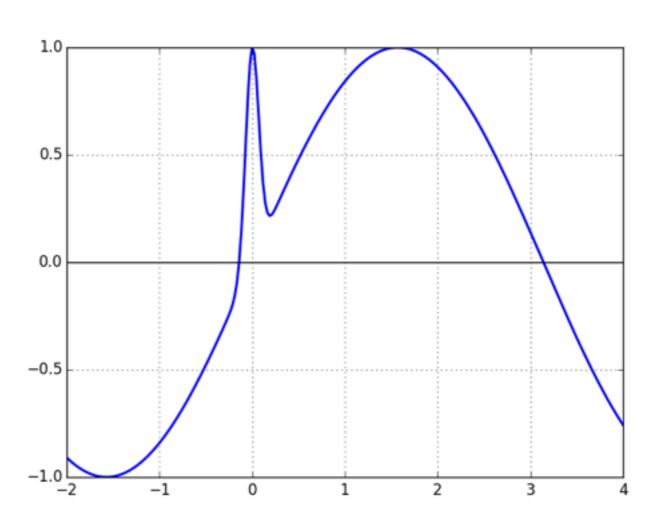
Recursion and Nested Threading

Adaptive Quadrature

• Problem: approximate

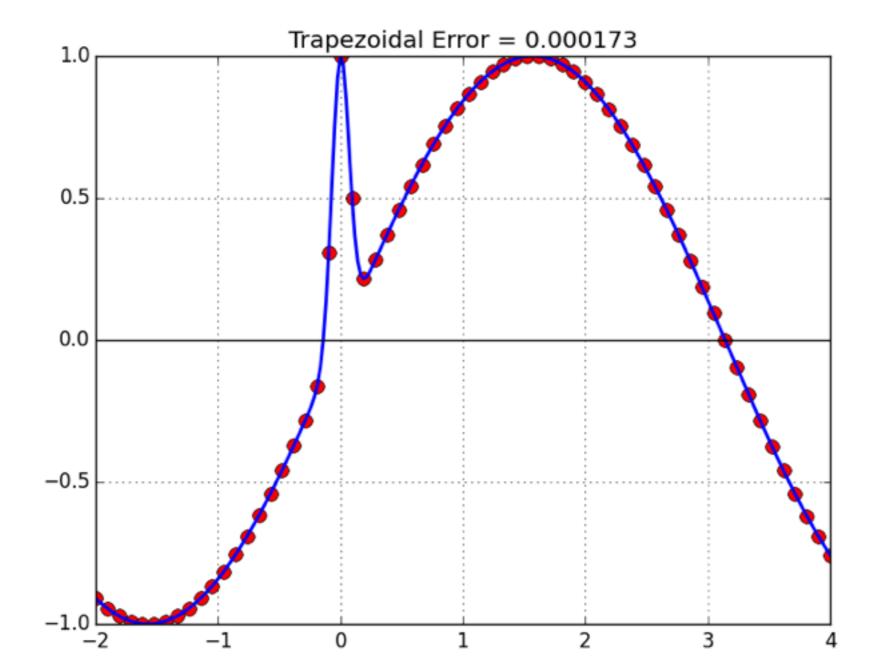
$$\int_{-2}^{4} e^{-\beta^{2}x^{2}} + \sin(x)dx = \left[\frac{\pi}{2\beta}\operatorname{erf}(\beta x) - \cos(x)\right]_{-2}^{4}$$

- where erf is the error function.
- beta = 10:



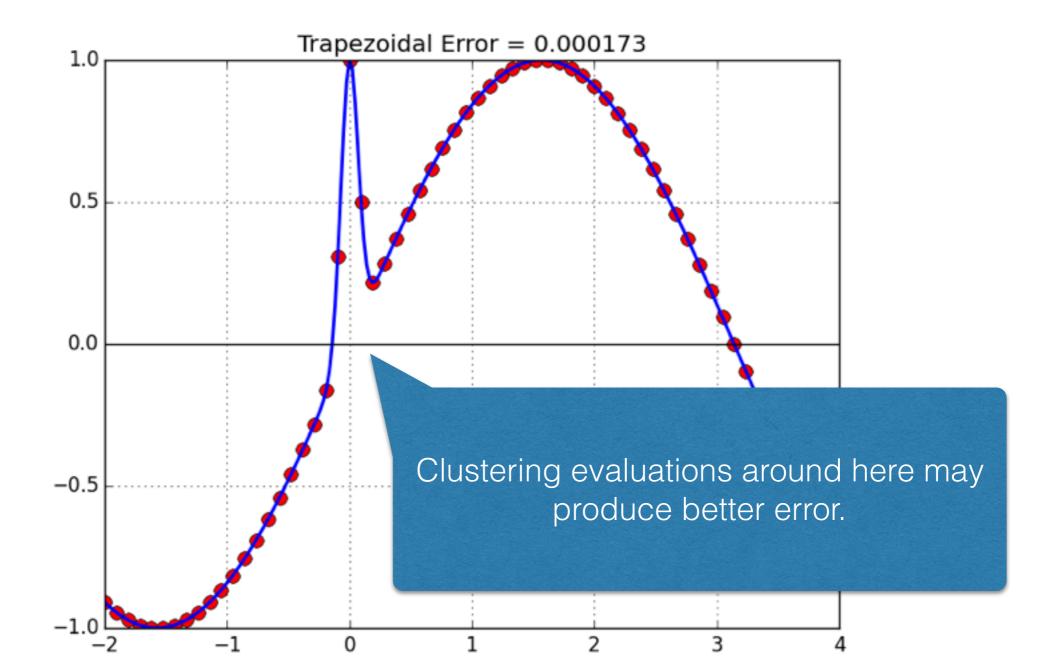
Estimating Integral

Trapezoidal rule: n=64 points, equally spaced



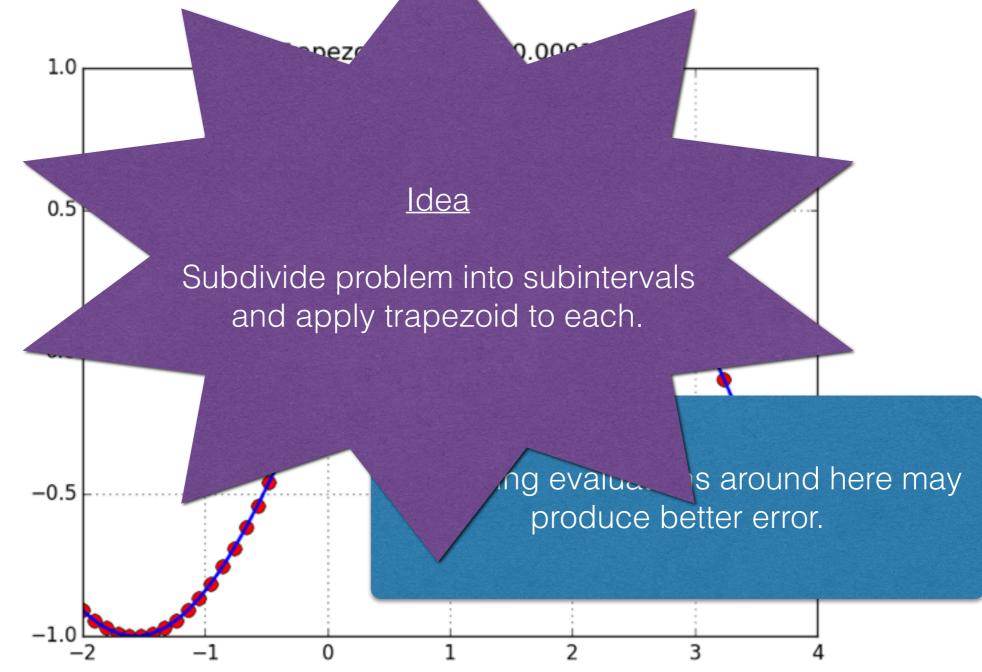
Estimating Integral

Trapezoidal rule: n=64 points, equally spaced



Estimating Integral

Trapezoidal rule: n=64 mints, equally spaced



Brief Aside

$$\int_{a}^{b} f(x)dx \approx \operatorname{trapz}(f, a, b)$$

- How to estimate approximation error?
- Compute both trapz and simps approx.
 - if abs. difference is large then bad approx.
 - if abs. difference is small then good approx.
 - Reason: error $\sim O(|b-a|^n)$

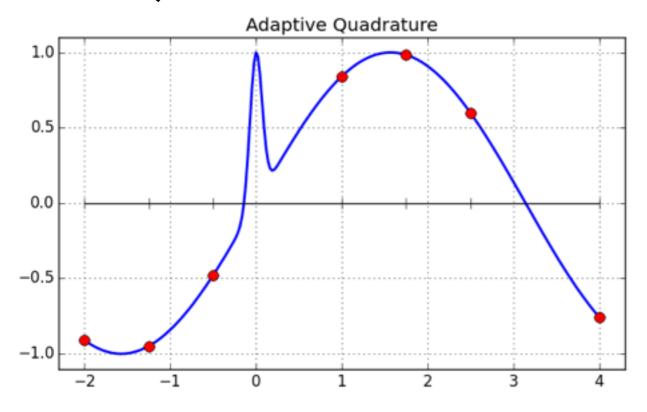
Central Idea

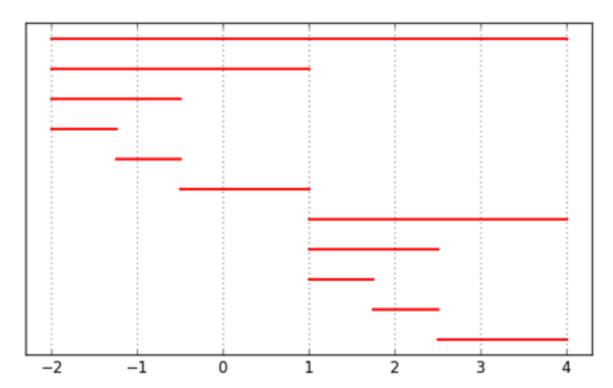
What if approx. error > tol?

$$\int_{a}^{b} f(x)dx = \int_{a}^{(a+b)/2} f(x)dx + \int_{(a+b)/2}^{b} f(x)dx$$

- Idea: split interval in half:
 - suppose error in each half is less than tol/2
 - then total error is less than tol
- If err > tol/2 then divide that half of the interval and repeat

Quadrature tol=0.5





integral: 0.198245

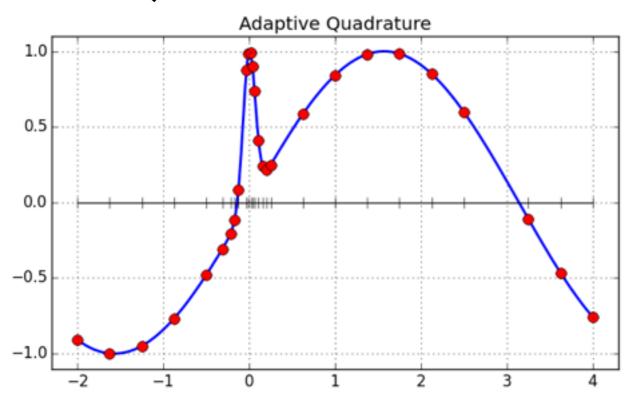
(actual: 0.414742)

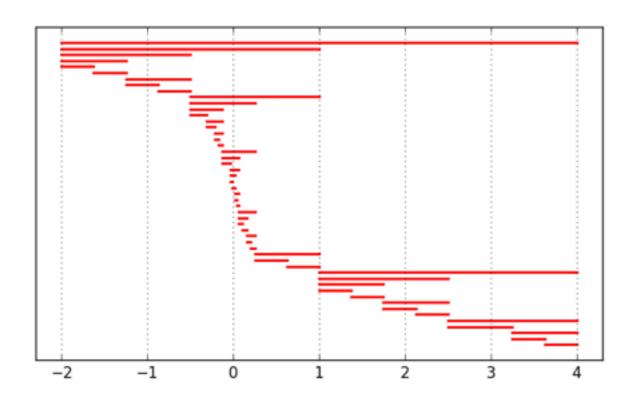
tolerance: 5.000000e-01

est error: 2.223337e-01

act error: 2.164973e-01

Quadrature tol=0.1





integral: 0.407417

(actual: 0.414742)

tolerance: 1.000000e-01

est error: 5.342968e-02

act error: 7.325371e-03

Demo

adaptive_serial.c and test_serial.c

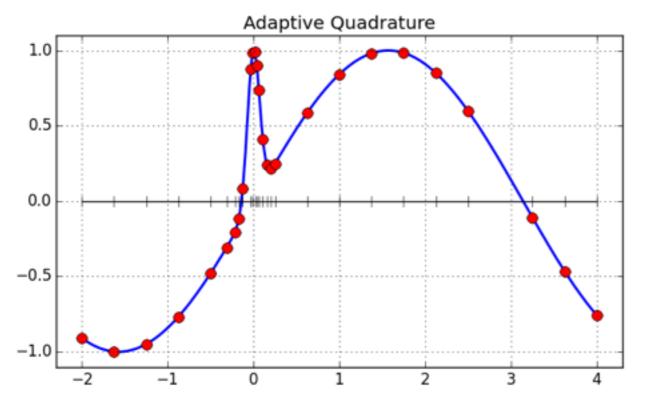
Parallelizing Using OpenMP

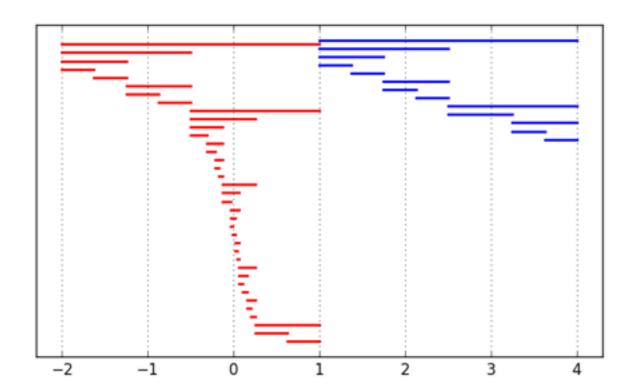
- Naive Idea: have each thread take a half of the domain
 - "domain decomposition pattern"
 - atomic / critical operations:
 - error calculation
 - integral summation

Demo

adaptive_serial.c and test_parallel.c

Par. Quad. tol=0.1





integral: 0.407417

(actual: 0.414742)

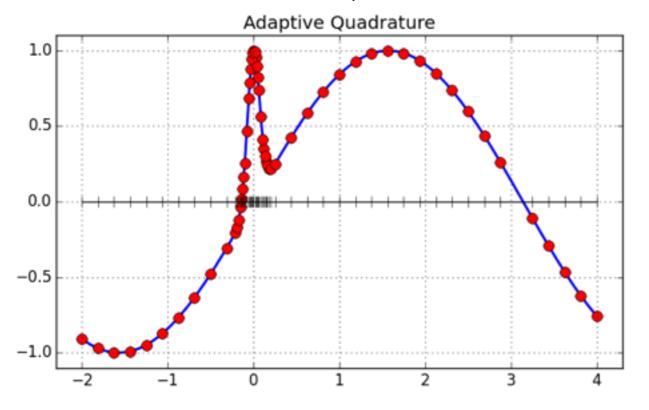
tolerance: 1.000000e-01

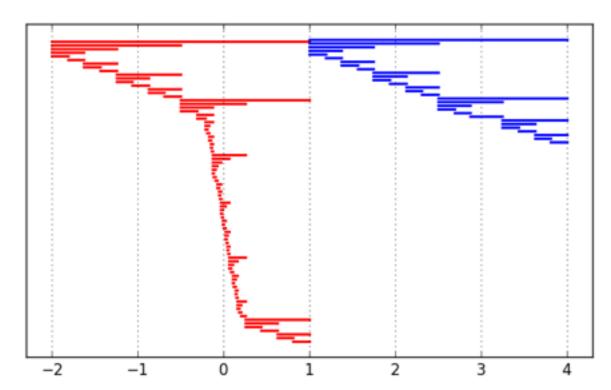
est error: 5.342968e-02

act error: 7.325371e-03

(Same as serial.)

Par. Quad. tol=0.02





integral: 0.414474

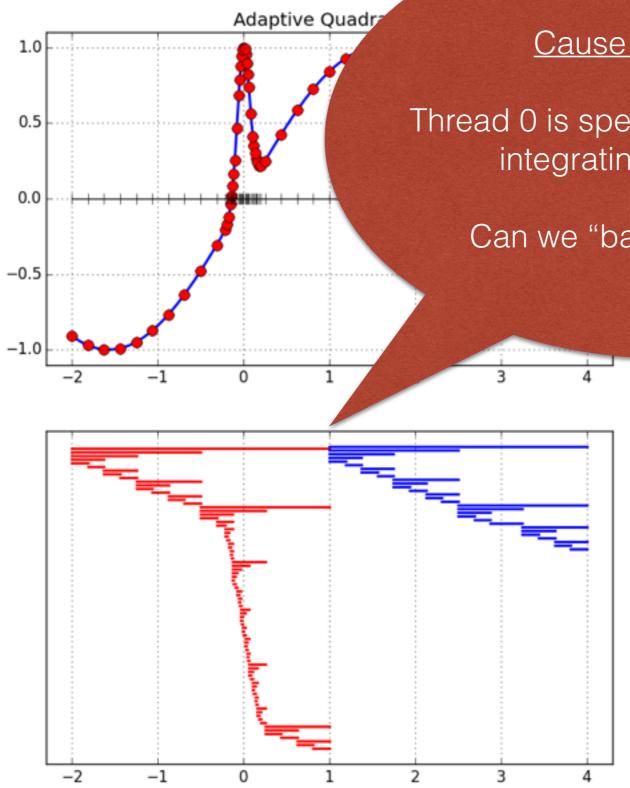
(actual: 0.414742)

tolerance: 2.000000e-02

est error: 1.189837e-02

act error: 2.678713e-04

Par. Quad 20.02



Cause for Concern

Thread 0 is spending almost 3x time integrating as Thread 1.

Can we "balance" the load?

0.414474

(actual: 0.414742)

tolerance: 2.000000e-02

est error: 1.189837e-02

act error: 2.678713e-04

 Work-sharing Construct (like omp for) — each "section" is given to next available thread

Work-sharing Construct (like omp for) — each "section" is given to next available thread

```
#pragma omp parallel sections
{
    #pragma omp section
    // block of code
    #pragma omp section
    // block of code
}
```

Allowed to combine parallel and sections into one.

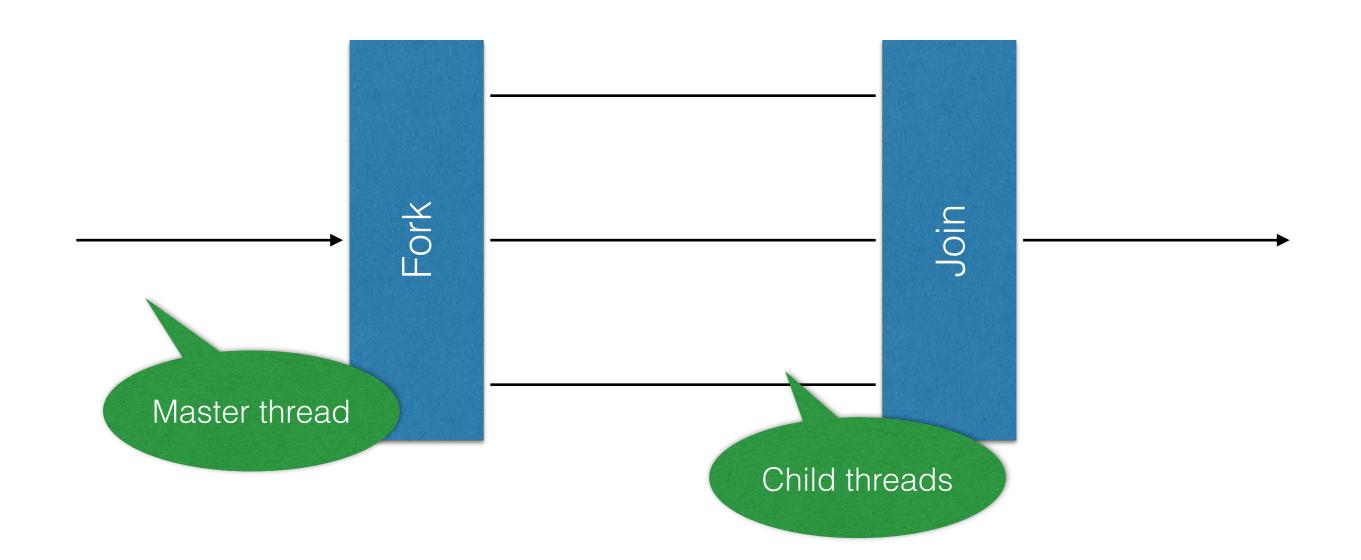
e.g. #pragma omp parallel for

- What happens if #threads =/= #sections?
 - If more threads than sections?

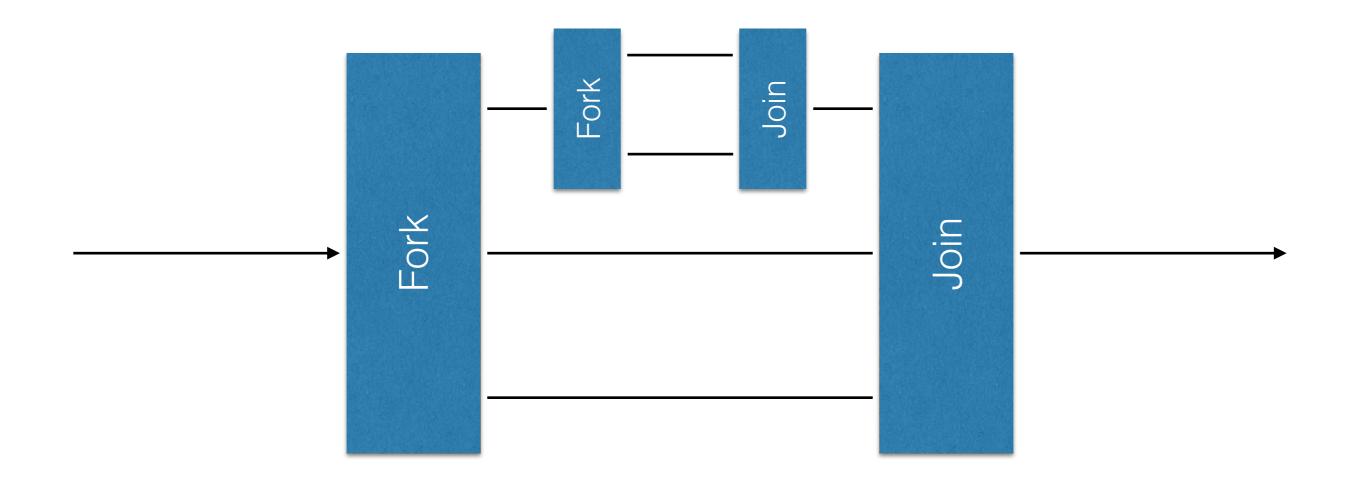
If more sections than threads?

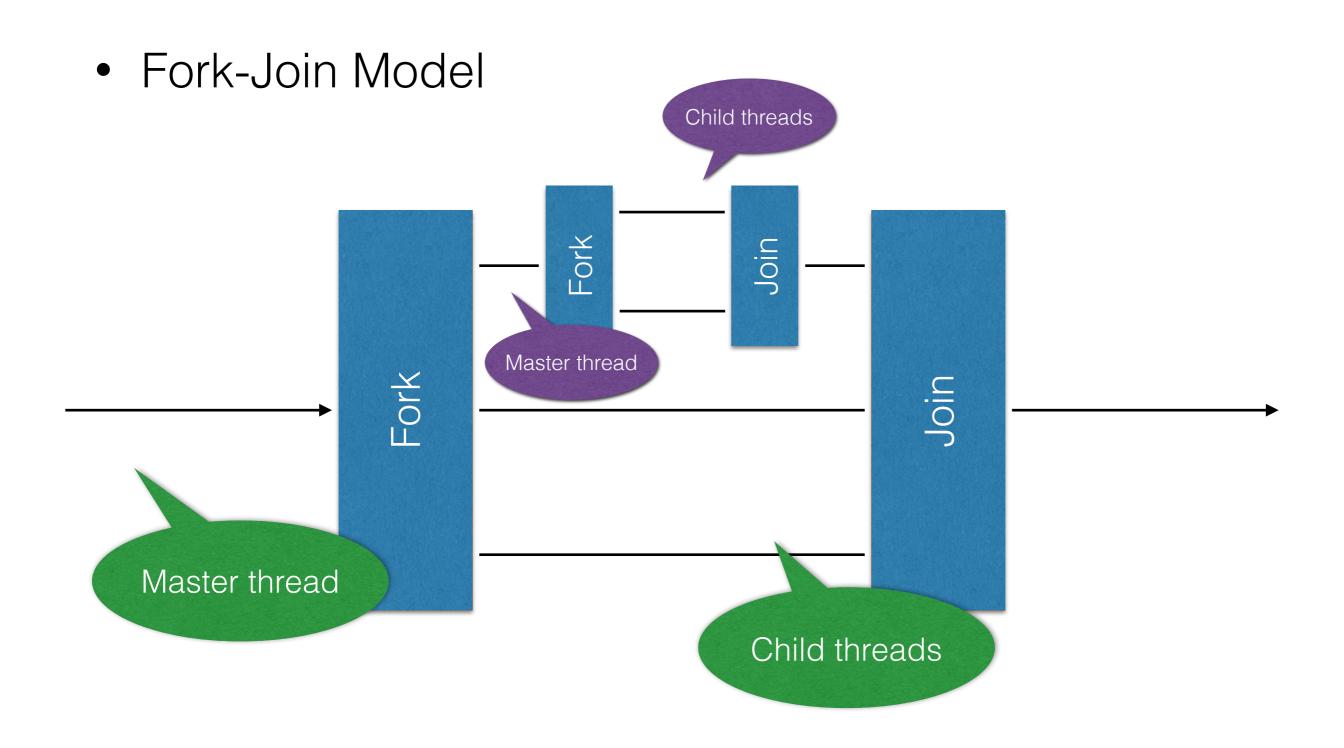
- What happens if #threads =/= #sections?
 - If more threads than sections?
 - Some threads will execute sections and some won't.
 - If more sections than threads?
 - OpenMP will allocate next available thread to next available section.

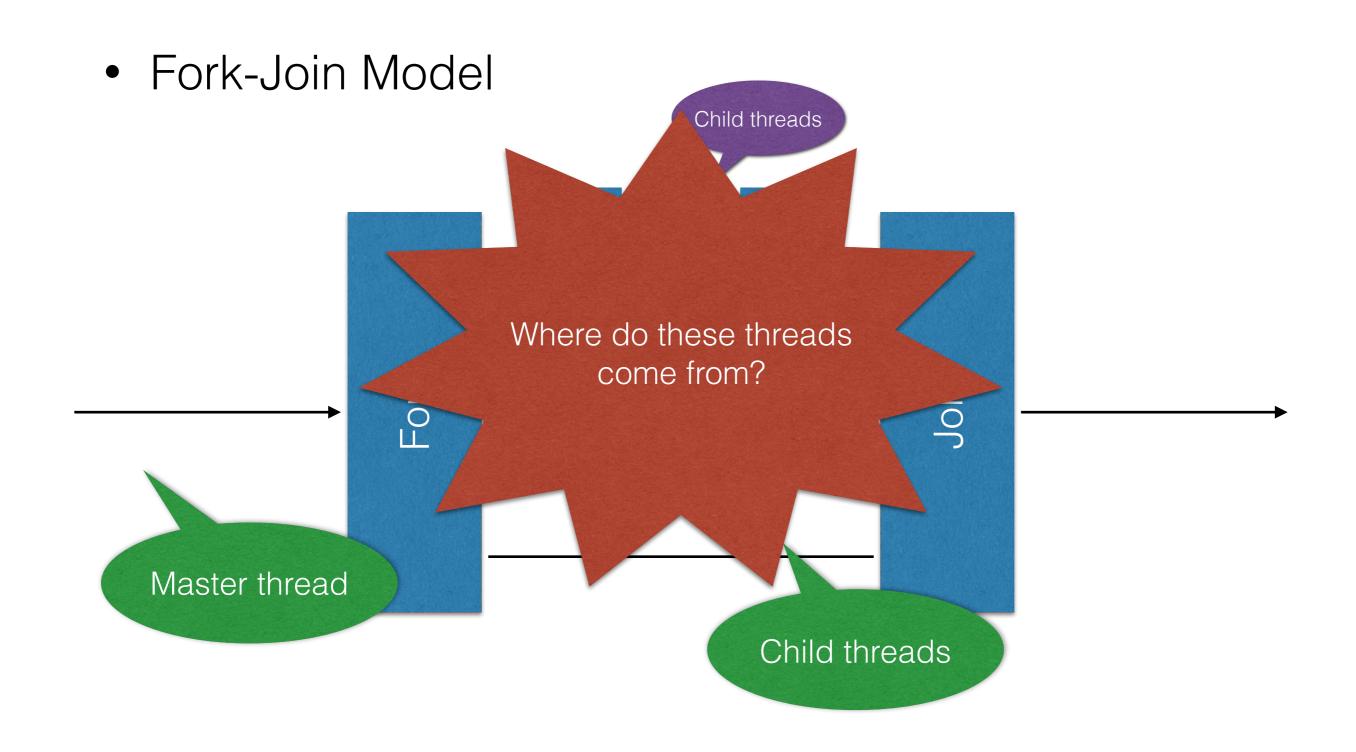
Fork-Join Model



Fork-Join Model







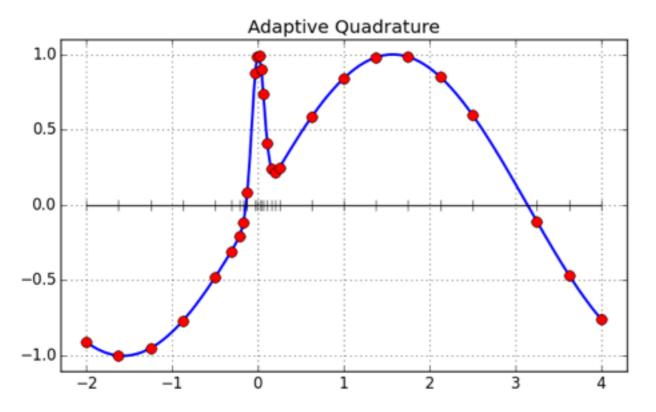
- OpenMP maintains a "pool" of threads
 - encounter omp parallel —> take idle threads
 - encounter another omp parallel —> take idle threads
- After JOIN, threads are returned to pool.

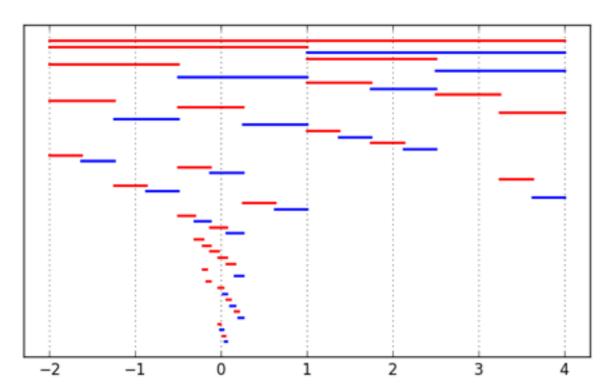
- Possible to create many total threads!
- Many ways to deal with this, see online resources.

Demo

adaptive_parallel.c and test_balanced.c

Balanced Quad. tol=0.1





 $num_threads = 2$

integral: 0.407417

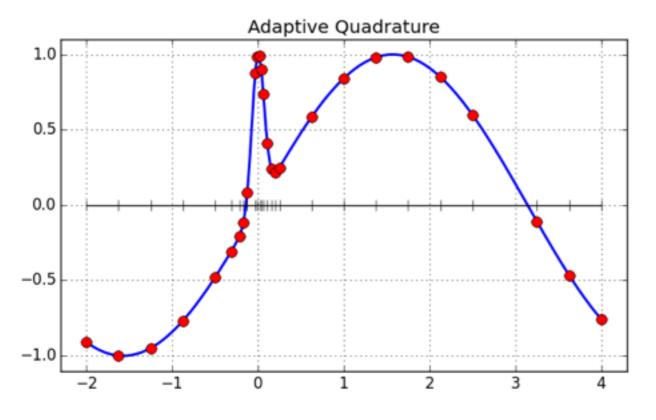
(actual: 0.414742)

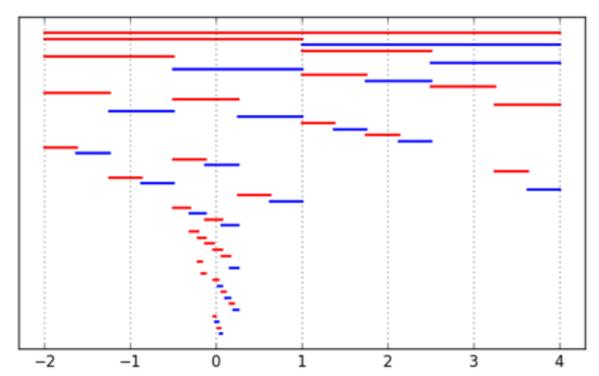
tolerance: 1.000000e-01

est error: 5.342968e-02

act error: 7.325371e-03

Balanced Quad. tol=0.1



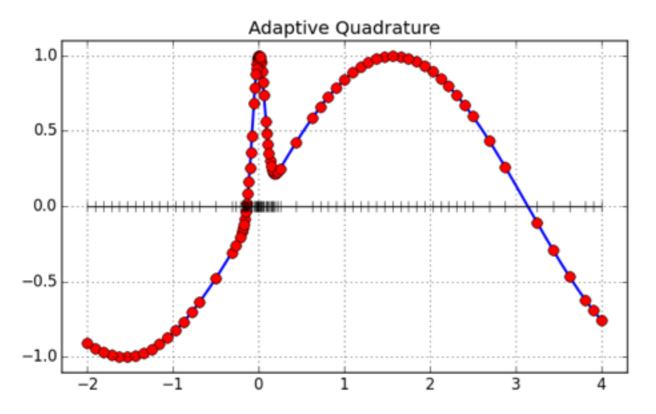


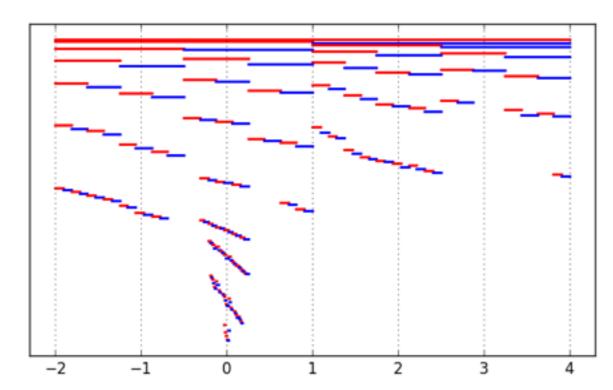
nu in (a to es

Work is shared evenly between Thread 0 and Thread 1.

(Each nested omp section given to next available thread.)

Balanced Quad. tol=0.01





 $num_threads = 2$

integral: 0.414743

(actual: 0.414742)

tolerance: 1.000000e-02

est error: 4.902247e-03

act error: 9.255080e-07

Thoughts on MPI

- OpenMP has built in tools for thread management
- Possible with MPI via MPI_Comm_spawn() and master —> child design pattern
- Many recursive functions can be re-implemented w/o recursion using a manual stack.
 - parallel producer consumer model

Thoughts on MPI

