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Lecture 13: Load balan

Previous lectures

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Several times in this module we have mentioned the concept of **load b**

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- Usually realised when **synchronisi**
- First encountered for the Mandelbrot set g
- Important for parallel performance for shared and distributed memory CPU, and GPU.

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Today's lecture

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Today we will look at load balancing more closely, and how to reduce

- <https://eduassistpro.github.io>^{tor, this}
- Understand how heterogeneity in the problem results in poor load balancing
- See how a **task scheduler** can improve **runtime**.
- Go through a concrete example of a **work pool**.

The Mandelbrot set (c.f. Lecture 3)

Code on Minerva: `Mandelbrot_MPI.c` plus makefile

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- Domain is $-2 \leq x \leq 2$ and $-2 \leq y \leq 2$.

- C

it

- P

the number of iterations.

- Here, the black region corresponds to a **high number of iterations**.

- **No upper bound** - some points will iterate **indefinitely** if allowed.

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Strip partitioning

Partition the domain [cf. last lecture] into horizontal strips¹.

Process 0



Process 3

Process p-1

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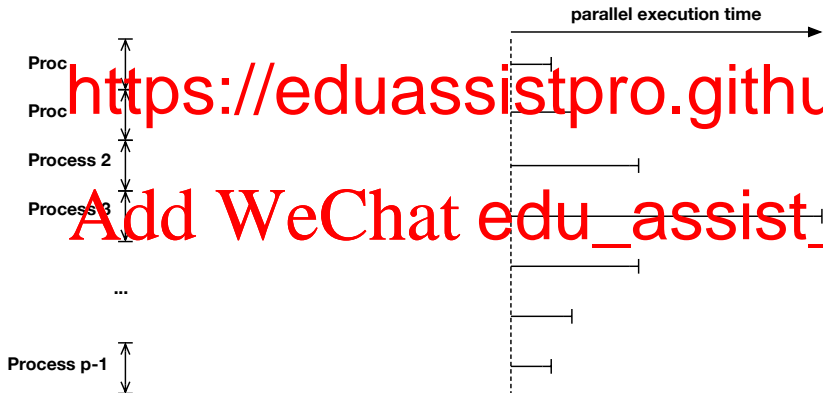
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¹Equivalent results for partitioning into vertical strips, or blocks.

Load imbalance

Because some pixels take longer to calculate the colour than others, the load is **unevenly distributed** across the processes:



Load balancing

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- Parallel execution time determined by the **last** processing unit to finish.

- r at least

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Definition

The goal of **load balancing** is for each **processing unit** (thread or process) to perform a **similar volume of computations**, and therefore finish at roughly the same time.

Up until now most problems we have encountered have been naturally load balanced

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For example, computing each p good

- Each unit performs n/p additions

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Note that the Mandelbrot set is a **map**

parallel problem (since there are no data dependencies).

- *Still* a challenge to attain good performance.

Static load balancing

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Definition

Som

a

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For the Mandelbrot set example, we could assign s to regions where the calculations should be



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Should improve load balancing.

However, an **exact** expression is not available. Therefore any such **heuristic** can only achieve **approximate** load balancing.

Static load balancing (*ideal case*)

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Process

Process

Process 2

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parallel execution time

d

d

saving compared to
unbalanced version

Dynamic load balancing

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Definition

D

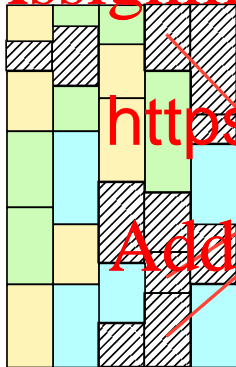
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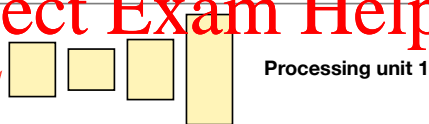
Basic idea:

- 1 Break the problem down into small
- 2 Each processing unit performs **on**
- 3 When it is complete it starts is assigned another task
- 4 repeat 3 until all tasks are complete

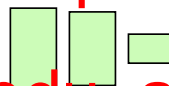
Full problem broken
down into tasks:



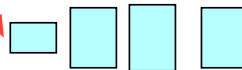
Tasks performed sequentially on
different threads/processes:



Processing unit 1



Processing unit 2



Processing unit 3

scheduler

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Functional or task parallelism

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Up to now, we have largely considered parallelising the same operation to a (large) data set.



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Now we are parallelising a number of tasks



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Called **task parallelism** or **functional parallelism**



Be warned that these terms are sometimes used to refer to slightly different concepts.



More on task parallelism in Lecture 19.

Work pools

The scheduler assigns tasks to processing units at runtime.



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To understand the role of a scheduler, we will look at a scheduler implemented in MPI - a **central**



One process (usually rank 0) performs the scheduling - the **main** process¹.



Remaining processes action the tasks - the **workers**¹.

¹You may see 'master' (for main) and 'slaves' (for workers) in the literature.

Worker pseudocode

Function workerProcess() in Mandelbrot_MPI.c

```
1 initialise(); // Including MPI_Init().
2
3 while( true )
4 {
5     // Wa
6     MP
7
8     // Is this a termination request?
9     if( message==TERMINATE ) break;
10
11     // Else perform calculation and send back to rank 0
12     result = actionTask( message );
13     MPI_Send( result, ... );
14 }
15
16 finalise(); // Including MPI_Finalize().
```

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Main process pseudocode (1)

Function `mainProcess()` in `Mandelbrot_MPI.c`

```
1 initialiseAndOpenWindow();  
2  
3 // Init  
4 int numA  
5  
6 // Send in  
7 for( p=1; p<numProcs; p++ )  
8 {  
9     MPI_Send(task, ..., p, ...);  
10    numActive++;  
11 }
```

For this Mandelbrot example, each task is a **row of pixel colours to be calculated**.

- Keep track with an incrementing variable `row`.

Main process pseudocode (2)

Function idle() in Mandelbrot_MPI.c

```
1 while( numActive > 0 )
2 {
3     // Get result from ANY worker process.
4     MP
5     nu
6
7     // Se
8     if( !finished )
9     {
10         MPI_Send( task, ..., status, MPI_SOURCE, ...
11         numActive++;
12     }
13
14     // Action the message.
15     actionResult( result );
16 }
```

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Main process pseudocode (3)

Function `idle()` in `Mandelbrot_MPI.c`

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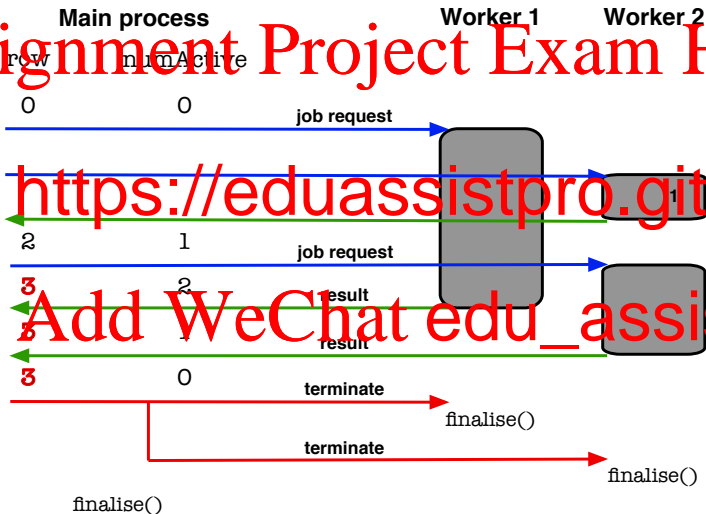
```
1 // Tell all workers to terminate.
2 for( p=1; p<numProcs; p++ )
3     MP
4
5 final
```

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- `MPI_ANY_SOURCE` in place of source in `recv` as a message from any process.
- Used `status.MPI_SOURCE` to recover process.
- Send next request **before** the (potentially slow) call to `actionResult()`.

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Example: 3 rows and 2 workers



Modern schedulers

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There are many more types of **work pool**, such as those with no 'main' process (*decentralised work pools*)¹.

A common

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- Performs tasks sequentially, starting from a queue
- Once the deque is empty, 'steals' a task from a randomly selected 'victim' (**work s**)

¹Wilkinson and Allen, *Parallel Programming* (Pearson, 2005).

OpenMP scheduler

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OpenMP can also schedule loops using its `schedule` clause:

```
1 #pragma omp parallel for schedule(dynamic, chunk)
2 for( i=0
```

This block is scheduled **at runtime**.

Can also

```
1 #pragma omp parallel for schedule(static, chunk)
```

There is also a **guided** option that decreases the chunk size exponentially **at runtime** to the final value

```
1 #pragma omp parallel for schedule(guided, chunk)
```

In all cases, `chunk` is optional and defaults to 1.

MIMD at last!

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Up until today we have mostly performed the **same** calculations on each processing unit.



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Today is the first clear¹ example where we
MIMD pattern **in software**.



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The **main** process perform entirely diff
workers (*division of labour*).

¹Ignoring trivial cases like e.g. rank 0 distributing global arrays.

Summary of distributed memory systems

Lec.	Content	Key points
8	Architectures and MPI	Clusters and supercomputers; interconnect network; starting with MPI.
9	Point-to-point communication	
10	Distributed communication	collective communication in MPI.
11	Reduction	Binary trees; OpenMP and MPI.
12	Asynchronous communication	Non-blocking partitioning a
13	Load balancing	Task parallelism; schedulers; work pools.

Next lecture we start looking at programming **general purpose graphics processing units** or GPGPUs.