Parallelism

COS 326
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Parallelism

What is it?

Why now? Today's technology trends

Some preliminary linguistic constructs
PARALLELISM:
WHAT IS IT?
Parallelism: What is it?

parallel computations involve doing many things at the same time instead of sequentially (one-after-the-other)
Flavors of Parallelism

Data Parallelism

– same computation being performed on a *collection* of independent items
– e.g., adding two vectors of numbers

Task Parallelism

– different computations/programs running at the same time
– e.g., running web server and database

Pipeline Parallelism

– assembly line:

map f over all items

map g over all items
Parallelism: performs many tasks *simultaneously*

- **purpose:** improves throughput (total jobs processed in allotted time)
- **mechanism:**
  - many independent computing devices
  - decrease run time of program by utilizing multiple cores or computers
- **eg:** running your web crawler on a cluster versus one machine.
Parallelism vs. Concurrency

**Parallelism**: performs many tasks *simultaneously*

- **purpose**: improves throughput (total jobs processed in allotted time)
- **mechanism**:
  - many independent computing devices
  - decrease run time of program by utilizing multiple cores or computers
- **eg**: running your web crawler on a cluster versus one machine.

**Concurrency**: mediates multi-party access to shared resources

- **purpose**: decrease response time
- **mechanism**:
  - switch between different threads of control
  - work on one thread when it can make useful progress; when it can't, suspend it and work on another thread
- **eg**: running your clock, editor, chat at the same time on a single CPU.
  - OS gives each of these programs a small time-slice (~10msec)
  - often *slows* throughput due to cost of switching contexts
- **eg**: don't block while waiting for I/O device to respond, but let another thread do useful CPU computation
Parallelism vs. Concurrency

**Parallelism:**
perform several independent tasks simultaneously

**Concurrency:**
mediate/multiplex access to shared resource

Many efficient programs use some parallelism and some concurrency.
PARALLELISM: WHY NOW?
UNDERSTANDING TECHNOLOGY TRENDS
Moore's Law

• Moore's Law: *The number of transistors you can put on a computer chip doubles (approximately) every couple of years.*

• Consequence for most of the history of computing: *All programs double in speed every couple of years.*

  – Why? Hardware designers are wicked smart.
  – They have been able to use those extra transistors to (for example) double the number of instructions executed per time unit, thereby processing speed of programs

• Consequence for application writers:

  – *watch TV for a while and your programs optimize themselves!*
  – new applications thought impossible became possible because of increased computational power
CPU Clock Speeds from 1993-2005
Next year’s machine is twice as fast!
CPU Clock Speeds from 1993-2005

Oops!
Power Dissipation

![Graph showing the increase in power density over time with various CPU models.](image)
Power Dissipation

(Dave-drawn curve.)
Moore’s Law still holds, so far, for transistors-per-chip.

But:
- clock speed,
- power
- performance/clock cycle all level off.

What do we do with all those transistors?
THE SOLUTION
Multi-core Hardware

- Core
  - ALU
  - ALU
- L1 cache
- L2 cache
- Main memory
GPUs

- There's nothing like video gaming to drive progress in computation!
- GPUs can have hundreds or even thousands of cores
- Three of the 5 most powerful supercomputers in the world take advantage of GPU acceleration.
- Scientists use GPUs for simulation and modelling
  - eg: protein folding and fluid dynamics
So...

Instead of trying to make your CPU go faster, Intel’s just going to pack more CPUs onto a chip.

– a few years ago: dual core (2 CPUs).
– a little more recently: 4, 6, 8 cores.
– Soon we may have hundreds or thousands on a chip.

In fact, that’s already happening with graphics chips.

– really good at simple data parallelism (many deep pipes)
– but they are *much* dumber than an Intel core.
– and right now, chew up a *lot* of power.
– watch for GPUs to get “smarter” and more power efficient, while CPUs become more like GPUs.
STILL MORE PROCESSORS: THE DATA CENTER
... so I thought to myself, why clutter your brain up with memories when I can store them in the cloud?
Data Centers: *Lots* of Connected Computers!
10s or 100s of thousands of computers connected together

• Motivated by new applications and scalable web services:
  – let's catalogue all N billion webpages in the world
  – let's all allow anyone in the world to search for the page he or she needs
  – let's process that search in less than a second

It's rather amazing. We are living science fiction.
Data Centers: Lots of Connected Computers

Computer containers for plug-and-play parallelism:
Sounds Great!

So my old programs will run 2x, 4x, 48x, 256x, 1024x faster?
Sounds Great!

So my old programs will run 2x, 4x, 48x, 256x, 1024x faster?
   – no way!
Sounds Great!

So my old programs will run 2x, 4x, 48x, 256x, 1024x faster?

– no way!

– to upgrade from Intel 386 to 486, the app writer and compiler writer did not have to do anything (much)
  • IA 486 interpreted the same sequential stream of instructions; it just did it faster
  • this is why we could watch TV while Intel engineers optimized our programs for us

– to upgrade from Intel 486 to dual core, we need to figure out how to split a single stream of instructions in to two streams of instructions that collaborate to complete the same task.
  • without work & thought, our programs don't get any faster at all
  • it takes ingenuity to generate efficient parallel algorithms from sequential ones
In Part: Functional Programming!

- Hadoop
- Naiad
- Pig
- Spark

Lightning-Fast Cluster Computing
PARALLEL AND CONCURRENT PROGRAMMING
• **Speedup**: the ratio of sequential program execution time to parallel execution time.

• If $T(p)$ is the time it takes to run a computation on $p$ processors

\[
\text{speedup}(p) = \frac{T(1)}{T(p)}
\]

• A parallel program has **perfect speedup** (aka **linear speedup**) if

\[
\frac{T(1)}{T(p)} = \text{speedup} = p
\]

• **Bad news**: Not every program can be effectively parallelized.
  – in fact, very few programs will scale with perfect speedups.
  – we certainly can't achieve perfect speedups automatically
  – limited by sequential portions, data transfer costs, ...
Most, but not all, parallel and concurrent programming models are far harder to work with than sequential ones:

- **They introduce nondeterminism**
  - the root of (almost all) evil
  - program parts suddenly have many different outcomes
    - they have different outcomes on different runs
    - debugging requires considering all of the possible outcomes
    - horrible heisenbugs hard to track down
  
- **They are nonmodular**
  - module A implicitly influences the outcomes of module B

- **They introduce new classes of errors**
  - race conditions, deadlocks

- **They introduce new performance/scalability problems**
  - busy-waiting, sequentialization, contention
Informal Error Rate Chart

regularity with which you shoot yourself in the foot
Informal Error Rate Chart

- Regularity with which you shoot yourself in the foot
- Null pointers, paucity of types, inheritance
- Manual memory management
- Kitchen sink + manual memory
- Heaven on earth
- Unstructured parallel or concurrent programming
Solid Parallel Programming Requires

1. Good sequential programming skills.
   – all the things we’ve been talking about: use modules, types, ...

2. Deep knowledge of the application.

3. **Pick a correct-by-construction parallel programming model**
   – whenever possible, a parallel model with semantics that coincides with sequential semantics
     • whenever possible, reuse well-tested libraries that hide parallelism
   – whenever possible, a model that cuts down non-determinism
   – whenever possible, a model with fewer possible concurrency bugs
   – if bugs can arise, know and use safe programming patterns

4. Careful engineering to ensure scaling.
   – unfortunately, there is sometimes a tradeoff:
     • reduced nondeterminism can lead to reduced resource utilization
   – synchronization, communication costs may need optimization
OUR FIRST PARALLEL PROGRAMMING MODEL: THREADS
Concurrent Threads with Locks: the classic shoot-yourself-in-the-foot concurrent programming model

- all the classic error modes

Why Threads?

- almost all programming languages will have a threads library
  - OCaml in particular!
- you need to know where the pitfalls are
- the assembly language of concurrent programming paradigms
  - we’ll use threads to build several higher-level programming models
Threads: an abstraction of a processor.

– programmer (or compiler) decides that some work can be done in parallel with some other work, e.g.:

```ocaml
let _ = compute_big_thing() in
let y = compute_other_big_thing() in
...
```

– we *fork* a thread to run the computation in parallel, e.g.:

```ocaml
let t = Thread.create compute_big_thing () in
let y = compute_other_big_thing () in
...
```
```ocaml
define	let
t = Thread.create f () in
let y = g () in
...
```

**Intuition in Pictures**

Processor 1:
- **Time 1:** `Thread.create`
- **Time 2:** `execute g ()`
- **Time 3:** `...`

Processor 2:
- **Time 1:** `(* doing nothing *)`
- **Time 2:** `execute f ()`
- **Time 3:** `...`
Suppose you have 2 available cores and you fork 4 threads....

In a typical multi-threaded system,

- the operating system provides *the illusion* that there are an infinite number of processors.
  - not really: each thread consumes space, so if you fork too many threads the process will die.

- it *time-multiplexes* the threads across the available processors.
  - about every 10 msec, it stops the current thread on a processor, and switches to another thread.
  - so a thread is really a *virtual processor*. 
Unfortunately, even if your computer has 2, 4, 6, 8 cores, OCaml cannot exploit them. It multiplexes all threads over a single core.

Hence, OCaml provides concurrency, but not parallelism. Why? Because OCaml (like Python) has no parallel “runtime system” or garbage collector. Other functional languages (Haskell, F#, ...) do.

Fortunately, when thinking about program correctness, it doesn’t matter that OCaml is not parallel -- I will often pretend that it is.

You can hide I/O latency, do multiprocess programming or distribute tasks amongst multiple computers in OCaml.
How do we get back the result that \( t \) is computing?

```ocaml
Thread.create : ('a -> 'b) -> 'a -> Thread.t

let t = Thread.create f () in
let y = g () in
...
```
First Attempt

```ocaml
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in

match !r with
| Some v -> (* compute with v and y *)
| None -> ???
```

What’s wrong with this?
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in
let rec wait() =
    match !r with
    | Some v -> v
    | None -> wait()

let v = wait() in
(* compute with v and y *)
Two Problems

First, we are *busy-waiting*.

- consuming cpu without doing something useful.
- the processor could be either running a useful thread/program or power down.

```ocaml
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in

let rec wait() =
    match !r with
    | Some v -> v
    | None -> wait()

in
let v = wait() in
(* compute with v and y *)
```
Second, an operation like \( r := \text{Some} \, v \) may not be *atomic*.

- \( r := \text{Some} \, v \) requires us to copy the bytes of \( \text{Some} \, v \) into the \texttt{ref} \( r \)
- we might see part of the bytes (corresponding to \( \text{Some} \)) before we’ve written in the other parts (e.g., \( v \)).
- So the waiter might see the wrong value.

```ocaml
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in
let rec wait() =
    match !r with
    | Some v -> v
    | None -> wait()

let v = wait() in
(* compute with v and y *)
```
An Aside: Atomicity

Consider the following:

```plaintext
let inc(r:int ref) = r := (!r) + 1
```

and suppose two threads are incrementing the same ref r:

**Thread 1**
```
inc(r);
!r
```

**Thread 2**
```
inc(r);
!r
```

If r initially holds 0, then what will Thread 1 see when it reads r?
The problem is that we can’t see exactly what instructions the compiler might produce to execute the code.

It might look like this:

**Thread 1**

R1 := load(p);
R1 := R1 + 1;
store R1 into r
R1 := load(p)

**Thread 2**

R1 := load(p);
R1 := R1 + 1;
store R1 into p
R1 := load(p)
Atomicity

But a clever compiler might optimize this to:

**Thread 1**

\[ R1 := \text{load}(p); \]
\[ R1 := R1 + 1; \]
\[ \text{store } R1 \text{ into } p \]
\[ R1 := \text{load}(r) \]

**Thread 2**

\[ R1 := \text{load}(p); \]
\[ R1 := R1 + 1; \]
\[ \text{store } R1 \text{ into } p \]
\[ R1 := \text{load}(r) \]
Furthermore, we don’t know when the OS might interrupt one thread and run the other.

Thread 1
R1 := load(p);
R1 := R1 + 1;
store R1 into p
R1 := load(r)

Thread 2
R1 := load(p);
R1 := R1 + 1;
store R1 into p
R1 := load(p)

(The situation is similar, but not quite the same on multi-processor systems.)
Atomicity

One possible interleaving of the instructions:

<table>
<thead>
<tr>
<th>Thread 1</th>
<th>Thread 2</th>
</tr>
</thead>
<tbody>
<tr>
<td>R1 := load(p);</td>
<td>R1 := load(p);</td>
</tr>
<tr>
<td>R1 := R1 + 1;</td>
<td>R1 := R1 + 1;</td>
</tr>
<tr>
<td>store R1 into r</td>
<td>store R1 into p</td>
</tr>
<tr>
<td>R1 := load(p)</td>
<td>R1 := load(p)</td>
</tr>
</tbody>
</table>

What answer do we get?
Another possible interleaving:

**Thread 1**
- \( R1 := \text{load}(p); \)
- \( R1 := R1 + 1; \)
- store \( R1 \) into \( p \)
- \( R1 := \text{load}(p) \)

**Thread 2**
- \( R1 := \text{load}(p); \)
- \( R1 := R1 + 1; \)
- store \( R1 \) into \( p \)
- \( R1 := \text{load}(p) \)

What answer do we get this time?
Another possible interleaving:

Thread 1
R1 := load(p);
R1 := R1 + 1;
store R1 into p
R1 := load(r)

Thread 2
R1 := load(p);
R1 := R1 + 1;
store R1 into p
R1 := load(p)

What answer do we get this time?

**Moral:** The system is responsible for *scheduling* execution of instructions.

**Moral:** This can lead to an enormous degree of *nondeterminism*. 
Even Worse ...

In fact, today’s multicore processors don’t treat memory in a *sequentially consistent* fashion. That means that we can’t even assume that what we will see corresponds to some interleaving of the threads’ instructions!

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**Beyond the scope of this class!** But the take-away is this: It’s not a good idea to use ordinary loads/stores to synchronize threads; you should use explicit synchronization primitives so the hardware and optimizing compiler don’t optimize them away.

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When Core1 stores to “memory”, it *lazily* propagates to Core2’s L1 cache. The load at Core2 might not see it, unless there is an explicit synchronization.
Even Worse

In fact, today’s multicore processors don’t treat memory in a *sequentially consistent* fashion. That means that *we can’t even assume that what we will see corresponds to some interleaving of the threads’ instructions!*

<table>
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**Beyond the scope of this class!** But the take-away is this: It’s not a good idea to use ordinary loads/stores to synchronize threads; you should use explicit synchronization primitives so the hardware and optimizing compiler don’t optimize them away.
The Happens Before Relation

We assume OCaml obeys a particular *Happens Before* relation:

**Rule 1:** Given two expressions (or instructions) in sequence: $e_1; e_2$ we know that $e_1$ happens before $e_2$.

**Rule 2:** Given a program:

```ocaml
let t = Thread.create f x in
....
Thread.join t;
e
```

we know that $(f \ x)$ happens before $e$. 
In Pictures

We know that for each thread the previous instructions must happen before the later instructions.

So for instance, inst\(_{1,1}\) must happen before inst\(_{1,2}\).
Thread 1
\[ t = \text{create f x} \]
\[ \text{inst}_{1,1}; \]
\[ \text{inst}_{1,2}; \]
\[ \text{inst}_{1,3}; \]
\[ \text{inst}_{1,4}; \]
\ldots
\[ \text{inst}_{1,n-1}; \]
\[ \text{inst}_{1,n}; \]
\[ \text{join t} \]

Thread 2
\[ \text{inst}_{2,1}; \]
\[ \text{inst}_{2,2}; \]
\[ \text{inst}_{2,3}; \]
\ldots
\[ \text{inst}_{2,m}; \]

We also know that the fork must happen before the first instruction of the second thread.
Thread 1
\[ t = \text{create f x} \]
\[ \text{inst}_{1,1}; \]
\[ \text{inst}_{1,2}; \]
\[ \text{inst}_{1,3}; \]
\[ \text{inst}_{1,4}; \]
... 
\[ \text{inst}_{1,n-1}; \]
\[ \text{inst}_{1,n}; \]
join t

Thread 2
\[ \text{inst}_{2,1}; \]
\[ \text{inst}_{2,2}; \]
\[ \text{inst}_{2,3}; \]
... 
\[ \text{inst}_{2,m}; \]

We also know that the fork must happen before the first instruction of the second thread.

And thanks to the join, we know that all of the instructions of the second thread must be completed before the join finishes.
However, in general, we do not know whether $\text{inst}_{1,i}$ executes before or after $\text{inst}_{2,j}$.

In general, \textit{synchronization instructions like fork and join reduce the number of possible interleavings}.

\textit{Synchronization cuts down nondeterminism}.

In the absence of synchronization we don’t know anything...
Calculate possible outcomes for a program by considering all of the possible interleavings of the *atomic* actions performed by each thread.

– Subject to the *happens-before* relation.
  
  • can’t have a child thread’s actions happening before a parent forks it.
  • can’t have later instructions execute earlier *in the same thread*.

– Here, *atomic* means indivisible actions.
  
  • For example, on most machines reading or writing a 32-bit word is atomic.
  • But, writing a multi-word object is usually *not* atomic.
  • Most operations like “b := b - w” are implemented in terms of a series of simpler operations such as
    
    – r1 = read(b); r2 = read(w); r3 = r1 − r2; write(b, r3)

Reasoning about all interleavings is *hard*. *just about impossible* for people

– Number of interleavings grows exponentially with number of statements.
– It can be hard to tell what is and isn’t atomic in a high-level language.
– **YOU ARE DOOMED TO FAIL IF YOU HAVE TO WORRY ABOUT THIS STUFF!**
Another approach to the coordination Problem

```
Thread.create : ('a -> 'b) -> 'a -> Thread.t

let t = Thread.create f () in
let y = g () in
...
```

How do we get back the result that \( t \) is computing?
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in
Thread.join t ;
match !r with
| Some v -> (* compute with v and y *)
| None -> failwith "impossible"
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in
Thread.join t ;
match !r with
| Some v -> (* compute with v and y *)
| None -> failwith "impossible"

Thread.join t causes the current thread to wait until the thread t terminates.
let r = ref None
let t = Thread.create (fun _ -> r := Some(f ())) in
let y = g() in
Thread.join t ;
match !r with
| Some v -> (* compute with v and y *)
| _ -> failwith "impossible"

So after the join, we know that any of the operations of t have **completed**.
FUTURES: A PARALLEL PROGRAMMING ABSTRACTION
The fork-join pattern we just saw is so common, we’ll create an abstraction for it:

```ocaml
module type FUTURE = 
sig
  type 'a future

  (* future f x forks a thread to run f(x) and stores the result in a future when complete *)
  val future : ('a->'b) -> 'a -> 'b future

  (* force f causes us to wait until the thread computing the future value is done and then returns its value. *)
  val force : 'a future -> 'a
end
```
Does that interface looks familiar .... ?
module Future : FUTURE =
struct
  type 'a future = {tid : Thread.t ;
                   value : 'a option ref }
end
module Future : FUTURE =
struct
    type 'a future = {tid : Thread.t ;
        value : 'a option ref }

    let future(f:'a->'b)(x:'a) : 'b future =
        let r = ref None in
        let t = Thread.create (fun () -> r := Some(f x)) ()
        in
        {tid=t ; value=r}
end
module Future : FUTURE =
struct
    type 'a future = {tid : Thread.t ;
        value : 'a option ref }

    let future(f:'a->'b)(x:'a) : 'b future =
        let r = ref None in
        let t = Thread.create (fun () -> r := Some(f x)) ()
        in
        {tid=t ; value=r}

    let force (f:'a future) : 'a =
        Thread.join f.tid ;
        match !(f.value) with
            | Some v -> v
            | None -> failwith "impossible!"
end
Now using Futures

```ocaml
let x = future f () in
let y = g () in
let v = force x in

(* compute with v and y *)
```
Back to the Futures

```ocaml
module type FUTURE =
  sig
    type 'a future
    val future : ('a->'b) -> 'a -> 'b future
    val force : 'a future -> 'a
  end

with futures library:

let x = future f () in
let y = g () in
let v = force x in
y + v

without futures library:

let r = ref None
let t = Thread.create
      (fun _ -> r := Some(f ()))
      ()
in
let y = g() in
Thread.join t ;
match !r with
  Some v -> y + v
| None -> failwith "impossible"

val f : unit -> int
val g : unit -> int
```

```ocaml
module type FUTURE =
  sig
    type 'a future
    val future : ('a->'b) -> 'a -> 'b future
    val force : 'a future -> 'a
  end

val f : unit -> int
val g : unit -> int
```
with futures library:

```ocaml
let x = future f () in
let y = g () in
let v = force x in
y + v
```

without futures library:

```ocaml
let r = ref None
let t = Thread.create
  (fun _ -> r := Some(f ()))
  ()
in
let y = g() in
Thread.join t ;
match !r with
  Some v -> y + v
| None -> failwith "impossible"
```

what happens if we delete these lines?
Back to the Futures

module type FUTURE =
sig
  type 'a future
  val future : ('a->'b) -> 'a -> 'b future
  val force : 'a future -> 'a
end

val f : unit -> int
val g : unit -> int

with futures library:

let x = future f () in
let y = g () in
let v = force x in
y + x

without futures library:

let r = ref None
let t = Thread.create
  (fun _ -> r := Some(f ()))
  ()
in
let y = g() in
Thread.join t ;
match !r with
  Some v -> y + v
| None -> failwith "impossible"

what happens if we use x and forget to force?
module type FUTURE =
  sig
    type 'a future

    val future : ('a->'b) -> 'a -> 'b future
    val force : 'a future -> 'a
  end

val f : unit -> int
val g : unit -> int

with futures library:

let x = future f () in
let y = g () in
let v = force x in
y + x

without futures library:

let r = ref None
let t = Thread.create
   (fun _ -> r := Some(f ()))
   ()
in
let y = g() in
Thread.join t ;
match !r with
  Some v -> y + v
| None -> failwith "impossible"

Moral: Futures + typing ensure entire categories of errors can’t happen -- you protect yourself from your own stupidity
module type FUTURE =
sig
  type 'a future

val future : ('a->'b) -> 'a -> 'b future
val force : 'a future -> 'a
end

with futures library:

let x = future f () in
let v = force x in
let y = g () in
y + x

what happens if you relocate force, join?

without futures library:

let r = ref None
let t = Thread.create
  (fun _ -> r := Some(f ()))
  ()
in
Thread.join t ;
let y = g() in
match !r with
  Some v -> y + v
| None -> failwith "impossible"
Back to the Futures

```ocaml
module type FUTURE =
  sig
    type 'a future
    val future : ('a->'b) -> 'a -> 'b future
    val force :'a future -> 'a
  end

val f : unit -> int
val g : unit -> int

with futures library:

let x = future f () in
let v = force x in
let y = g () in
y + x

Moral: Futures are not a universal savior

without futures library:

let r = ref None
let t = Thread.create
  (fun _ -> r := Some(f ()))
  ()
in
Thread.join t ;
let y = g() in
match !r with
  Some v -> y + v
| None -> failwith "impossible"
```
An Example: Mergesort on Arrays

```ocaml
let mergesort (cmp:'a->'a->int)
    (arr : 'a array) : 'a array =
let rec msort (start:int) (len:int) : 'a array =
    match len with
    | 0 -> Array.of_list []
    | 1 -> Array.make 1 arr.(start)
    | _ -> let half = len / 2 in
       let a1 = msort start half in
       let a2 = msort (start + half)
            (len - half) in
        merge a1 a2

and merge (a1:'a array) (a2:'a array) : 'a array =
...
```
An Example: Mergesort on Arrays

```ocaml
let mergesort (cmp:'a->'a->int)
  (arr : 'a array) : 'a array =
let rec msort (start:int) (len:int) :
  match len with
  | 0 -> Array.of_list []
  | 1 -> Array.make 1 arr.(start)
  | _ -> let half = len / 2 in
    let a1 = msort start half in
    let a2 = msort (start + half) (len - half) in
    merge a1 a2

and merge (a1:'a array) (a2:'a array) : 'a array = ...
```

Opportunity for parallelization
let mergesort (cmp:'a->'a->int) (arr : 'a array) : 'a array =

let rec msort (start:int) (len:int) : 'a array =
    match len with
    | 0 -> Array.of_list []
    | 1 -> Array.make 1 arr.(start)
    | _ -> let half = len / 2 in
      let a1_f = Future.future (msort start) half in
      let a2 = msort (start + half)(len - half) in
      merge (Future.force a1_f) a2

and merge (a1:'a array) (a2:'a array) : 'a array =
Divide-and-Conquer

This is an instance of a basic *divide-and-conquer* pattern in parallel programming

– take the problem to be solved and divide it in half
– fork a thread to solve the first half
– simultaneously solve the second half
– synchronize with the thread we forked to get its results
– combine the two solution halves into a solution for the whole problem.

**Warning:** the fact that we only had to rewrite 2 lines of code for mergesort made the parallelization transformation look deceptively easy

– we also had to verify that any two threads did not touch overlapping portions of the array -- if they did we would have to again worry about scheduling nondeterminism
Caveats

There is some overhead for creating a thread.
   – On uniprocessor, parallel code \textit{slower} than sequential code.

Even on a multiprocessor, we do \textit{not always} want to fork.
   – when the subarray is small, faster to sort it sequentially than to fork
     • similar to using insertion sort when arrays are small vs. quicksort
   – this is known as a \textit{granularity problem}
     • more parallelism than we can effectively take advantage of.
In a good implementation of futures, a compiler and run-time system might look to see whether the cost of doing the fork is justified by the amount of work that will be done. Today, it’s up to you to figure this out… 😞

– typically, use parallel divide-and-conquer until:
  (a) we have generated \textit{at least} as many threads as there are processors
  – often \textit{more threads} than processors because different jobs take different amounts of time to complete and we would like to keep all processors busy
  (b) the sub-arrays have gotten small enough that it’s not worth forking.

We’re not going to worry about these performance-tuning details but rather focus on the distinctions between \textit{parallel} and \textit{sequential algorithms}. 
Another Example

```ocaml
type 'a tree = Leaf | Node of 'a node
and 'a node = {left : 'a tree ;
    value : 'a ;
    right : 'a tree }

let rec fold (f:'a -> 'b -> 'b -> 'b) (u:'b) (t:'a tree) : 'b =
    match t with
    | Leaf -> u
    | Node n ->
        f n.value (fold f u n.left) (fold f u n.right)

let sum (t:int tree) = fold (+) 0 t
```
type 'a tree = Leaf | Node of 'a node

and 'a node = {left : 'a tree ;
  value : 'a ;
  right : 'a tree }

let rec pfold (f:'a -> 'b -> 'b -> 'b) (u:'b) (t:'a tree) : 'b =

  match t with
  | Leaf -> u
  | Node n ->
    let l_f = Future.future (pfold f u) n.left in
    let r = pfold f u n.right in
    f n.value (Future.force l_f) r

let sum (t:int tree) = pfold (+) 0 t
If the tree is unbalanced, then we’re not going to get the same speedup as if it’s balanced.

Consider the degenerate case of a list.

– The forked child will terminate without doing any useful work.
– So the parent is going to have to do all that work.
– Pure overhead... 😞

In general, lists are a horrible data structure for parallelism.

– *we can’t cut the list in half in constant time*
– for arrays and trees, we can do that (assuming the tree is balanced.)
type 'a tree = Leaf | Node of 'a node
and 'a node = { left : 'a tree ;
value : 'a ;
right : 'a tree }

let rec pfold (f:'a -> 'b -> 'b -> 'b) (u:'b) (t:'a tree) : 'b =
  match t with
  | Leaf -> u
  | Node n ->
    let l_f = Future.future (pfold f u) n.left in
    let r = pfold f u n.right in
    f n.value (Future.force l_f) r

let print (t:int tree) =
  pfold (fun n _ _ -> Printf.print "%d\n" n) ()
If code is purely functional, then it never matters in what order it is run. If \( f() \) and \( g() \) are pure then all of the following are equivalent:

\[
\begin{align*}
\text{let } x &= f() \text{ in} \\
\text{let } y &= g() \text{ in} \\
& e
\end{align*}
\]

\[
\begin{align*}
\text{let } y &= g() \text{ in} \\
\text{let } x &= f() \text{ in} \\
& e
\end{align*}
\]

\[
\begin{align*}
\text{let } y\_g &= \text{future } g() \text{ in} \\
\text{let } x &= \text{force } x\_f \text{ in} \\
& e
\end{align*}
\]

\[
\begin{align*}
\text{let } x\_f &= \text{future } f() \text{ in} \\
\text{let } y &= \text{force } y\_g \text{ in} \\
& e
\end{align*}
\]

As soon as we introduce \textit{side-effects}, the order starts to matter.

- This is why, IMHO, \textit{imperative} languages where even the simplest of program phrases involves a side effect, are doomed.
- Of course, we’ve been saying this for 30 years!
- See J. Backus’s Turing Award lecture, \textit{“Can Programming be Liberated from the von Neumann Style? A Functional Style and Its Algebra of Programs.”} 
  
http://www.cs.cmu.edu/~crary/819-f09/Backus78.pdf
Reasoning about the correctness of pure parallel programs that include futures is easy -- no harder than ordinary, sequential programs. (Reasoning about their performance may be harder.)

Reasoning about shared variables and synchronization is hard in general, but futures are a discipline for getting it right.

Much of programming-language design is the art of finding good disciplines in which it’s harder* to write bad programs.

Even aside from PL design, the same is true of software engineering with Abstract Data Types: if you engineer disciplines into your interfaces, it is harder for the user to get it wrong.

*but somebody will always find a way...