#### 15-210: Parallelism in the Real World

- Types of paralellism
- Parallel Thinking
- Nested Parallelism
- Examples (Cilk, OpenMP, Java Fork/Join)
- Concurrency

# Cray-1 (1976): the world's most expensive love seat



# Data Center: Hundred's of thousands of computers







#### Since 2005: Multicore computers

#### AMD Opteron (sixteen-core) Model 6274

by AMD

★★★☆ < (1 customer review)</p>

List Price: \$693.00

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You Save: \$93.01 (13%)

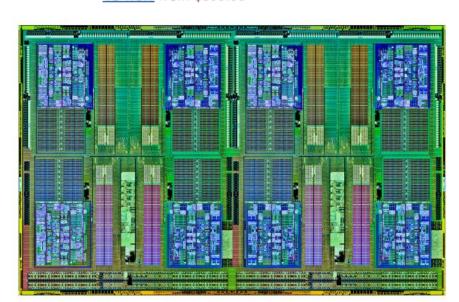


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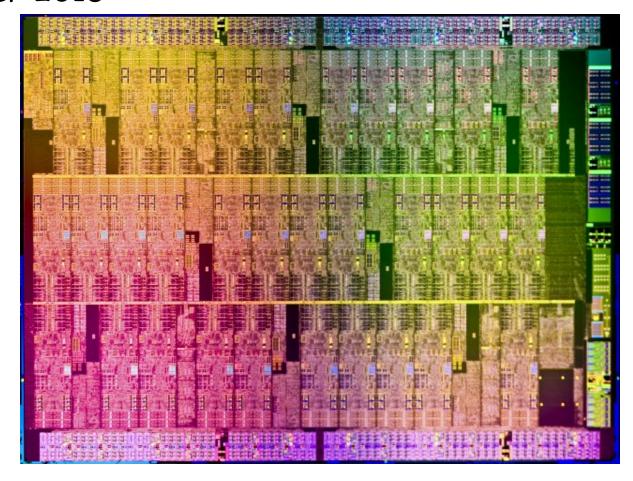
#### Xeon Phi: Knights Landing

72 Cores, x86, 500Gbytes of memory bandwidth Summer 2015



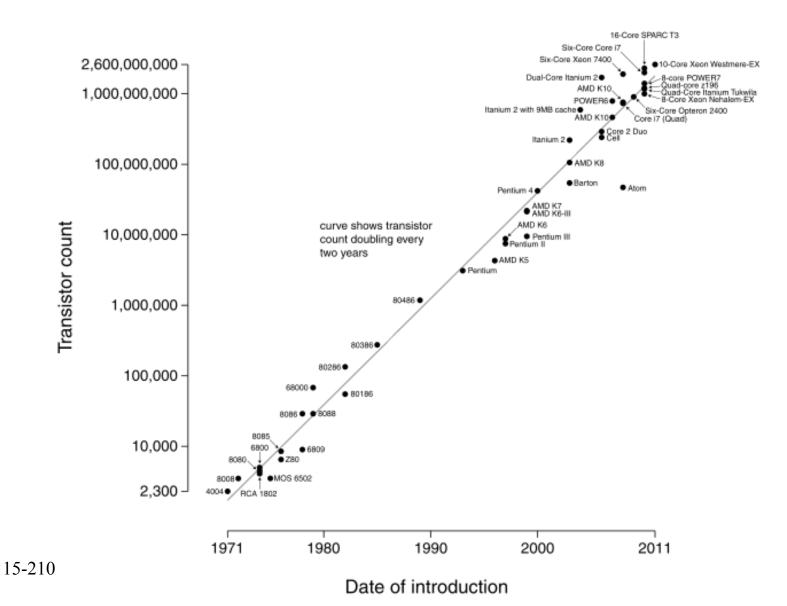
## Xeon Phi: Knights Landing

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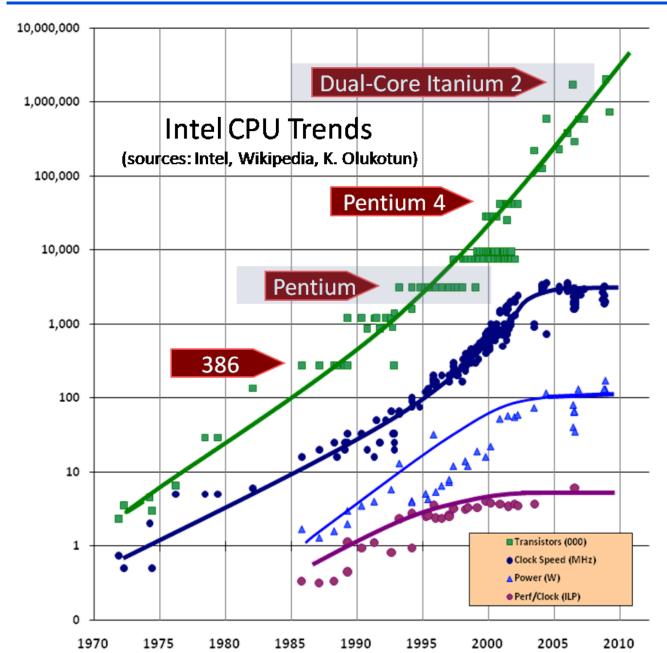
#### Moore's Law

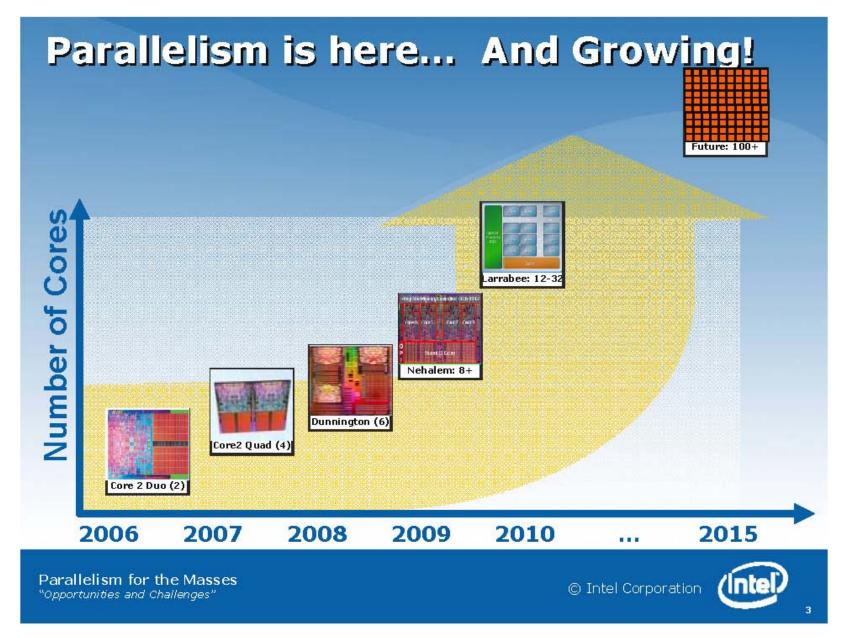
#### Microprocessor Transistor Counts 1971-2011 & Moore's Law



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#### Moore's Law and Performance





# 64 core blade servers (\$6K) (shared memory)

#### AMD Opteron (sixteen-core) Model 6274

by AMD

★★★☆ ▼ (1 customer review)

List Price: \$693.00

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 $\times 4 =$ 



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#### 1024 "cuda" cores





# EVGA GeForce GTX 590 Classified : 3DVI/Mini-Display Port SLI Ready Lii 03G-P3-1596-AR

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#### In Stock.

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Only 1 left in stock--order soon.

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# Samsung Galaxy S IV is now Official: Octa-Core CPU, 5" Full HD Display & 13MP Camera

Follow: Phones GT-I9500 Samsung Display Samsung Exynos Samsung Galaxy S IV Samsung Mobile Unpacked 2013



Samsung has just announced the Samsung Galaxy S4 at their Mobile Unpacked Event 2013 Episode 1 in New York, USA. The Galaxy S4 features a stunning 4.99" Full HD (1920×1080) SuperAMOED display. With a 441 ppi pixel density, your eyes won't be able to distinguish the pixels, which ensures excellent visual comfort. Even though the Galaxy S4 has a large display and a massive battery of 2,600 MAh, it's only 7.9mm thick. Samsung's latest

flagship device is PACKED with powerful components, consisting of Samsung's latest Exynos 5 Octa-Core (5410) CPU based on ARM's big.LITTLE technology with Quad Cortex-

# Intel Has a 48-Core Chip for Smartphones and Tablets

By Wolfgang Gruener OCTOBER 31, 2012 9:20 AM - Source: Computerworld

Intel has developed a prototype of a 48-core processor for smartphones. Before you ask: No, you can't buy a 48-core smartphone next year.



#### Parallel Hardware

#### Many forms of parallelism

- Supercomputers: large scale, shared memory
- Clusters and data centers: large-scale, distributed memory
- Multicores: tightly coupled, smaller scale
- GPUs, on chip vector units
- Instruction-level parallelism

Parallelism is important in the real world.

# Key Challenge: Software (How to Write Parallel Code?)

At a high-level, it is a two step process:

- Design a work-efficient, low-span parallel algorithm
- Implement it on the target hardware

In reality: each system required different code because programming systems are immature

- Huge effort to generate efficient parallel code.
  - Example: Quicksort in MPI is <u>1700 lines</u> of code, and about the same in CUDA
- Implement one parallel algorithm: a whole thesis.

Take 15-418 (Parallel Computer Architecture and Prog.)

## 15-210 Approach

Enable parallel thinking by raising abstraction level

- I. Parallel thinking: Applicable to many machine models and programming languages
- II. Reason about correctness and efficiency of algorithms and data structures.

## Parallel Thinking

Recognizing true dependences: unteach sequential programming.

Parallel algorithm-design techniques

- Operations on aggregates: map/reduce/scan
- Divide & conquer, contraction
- Viewing computation as DAG (based on dependences)

Cost model based on work and span

## Quicksort from Aho-Hopcroft-Ullman (1974)

procedure QUICKSORT(5):

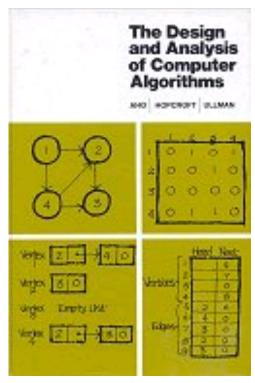
if S contains at most one element then return S

else

begin

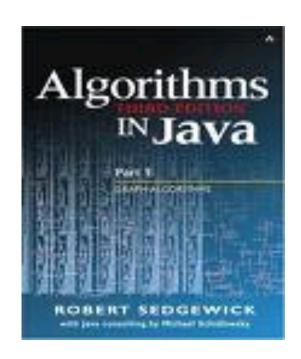
choose an element a randomly from S; let S<sub>1</sub>, S<sub>2</sub> and S<sub>3</sub> be the sequences of elements in S less than, equal to, and greater than a, respectively; return (QUICKSORT(S<sub>1</sub>) followed by S<sub>2</sub> followed by QUICKSORT(S<sub>3</sub>))

end



#### Quicksort from Sedgewick (2003)

```
public void quickSort(int[] a, int left, int right) {
    int i = left-1; int j = right;
    if (right <= left) return;</pre>
    while (true) {
      while (a[++i] < a[right]);
      while (a[right] < a[--j])
        if (j==left) break;
      if (i \ge j) break;
      swap(a,i,j); }
    swap(a, i, right);
    quickSort(a, left, i - 1);
    quickSort(a, i+1, right); }
```



## Styles of Parallel Programming

Data parallelism/Bulk Synchronous/SPMD Nested parallelism: what we covered Message passing Futures (other pipelined parallelism) General Concurrency

#### Nested Parallelism

# Nested Parallelism = arbitrary nesting of parallel loops + fork-join

- Assumes no synchronization among parallel tasks except at joint points.
- Deterministic if no race conditions

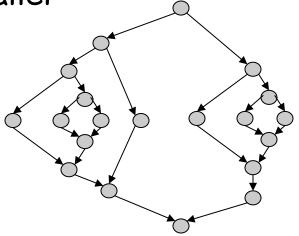
#### Advantages:

- Good schedulers are known
- Easy to understand, debug, and analyze costs
- Purely functional, or imperative...either works

#### Serial Parallel DAGS

Dependence graphs of nested parallel computations are

series parallel



Two tasks are parallel if not reachable from each other.

A data race occurs if two parallel tasks are involved in a race if they access the same location and at least one is a write.

#### Nested Parallelism: parallel loops

#### Nested Parallelism: fork-join

```
cobegin {
                              Dates back to the 60s. Used in
  S1;
                                 dialects of Algol, Pascal
  S2;}
                              Java fork-join framework
coinvoke(f1,f2)
                              Microsoft TPL (C#,F#)
Parallel.invoke(f1,f2)
#pragma omp sections
                              OpenMP (C++, C, Fortran, ...)
  #pragma omp section
  S1;
  #pragma omp section
  S2;
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                                                       Page24
```

## Nested Parallelism: fork-join

```
spawn S1;
S2;
                     cilk, cilk+
sync;
(exp1 | exp2)
                     Various functional
                       languages
plet
  x = exp1
                     Various dialects of
  y = exp2
                       ML and Lisp
in
  exp3
```

#### Cilk vs. what we've covered

ML:  $val(a,b) = par(fn() \Rightarrow f(x), fn() \Rightarrow g(y))$ 

Psuedocode: val (a,b) = (f(x) || g(y))

Cilk:  $cilk\_spawn a = f(x)$ ;

b = g(y);

cilk\_sync;

Fork Join

ML: S = map f A

Psuedocode:  $S = \langle f x : x \text{ in } A \rangle$ 

Cilk:  $cilk_{for}(int i = 0; i < n; i++)$ 

S[i] = f(A[i]) Page 26

#### Cilk vs. what we've covered

ML: S = tabulate f n

Psuedocode: S = <f i: i in <0,..n-1>> Tabulate

Cilk: cilk\_for (int i = 0; i < n; i++)

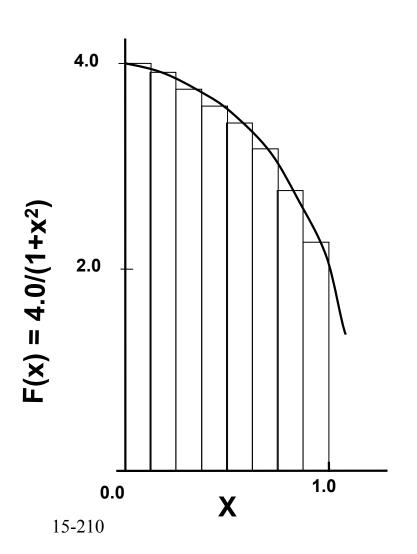
S[i] = f(i)

## Example Cilk

```
int fib (int n) {
  if (n<2) return (n);
  else {
    int x,y;
    x = cilk_spawn fib(n-1);
    y = cilk_spawn fib(n-2);
    cilk_sync;
    return (x+y);
  }
}</pre>
```

# Example OpenMP: Numerical Integration

Mathematically, we know that:



$$\int_{0}^{1} \frac{4.0}{(1+x^2)} dx = \pi$$

We can approximate the integral as a sum of rectangles:

$$\sum_{i=0}^{N} F(x_i) \Delta x \approx \pi$$

where each rectangle has width  $\Delta x$  and height  $F(x_i)$  at the middle of interval i.

#### The C code for Approximating PI

```
static long num_steps = 100000;
double step;
void main ()
       int i; double x, pi, sum = 0.0;
       step = 1.0/(double) num steps;
       x = 0.5 * step;
       for (i=0;i<= num_steps; i++){
              x+=step;
              sum += 4.0/(1.0+x*x);
       pi = step * sum;
```

## The C/openMP code for Approx. PI

```
#include <omp.h>
static long num_steps = 100000;
                                      double step;
void main ()
                                                 Private clause
    int i;
                double x, pi, sum = 0.0;
                                               creates data local to
    step = 1.0/(double) num_steps;
                                                   a thread
#pragma omp parallel for private(i, x) reduction(+:sum)
    for (i=0;i<= num steps; i++){
        x = (i+0.5)*step;
        sum = sum + 4.0/(1.0+x*x);
    pi = step * sum;
                                            Reduction used to
                                                manage
                                             dependencies
```

#### Example: Java Fork/Join

```
class Fib extends FJTask {
   volatile int result; // serves as arg and result
   int n;
   Fib(int n) { n = n; }
   public void run() {
      if (n \le 1) result = n;
      else if (n <= sequentialThreshold) number = seqFib(n);
      else {
         Fib f1 = \text{new Fib}(n - 1);
         Fib f2 = \text{new Fib}(n - 2);
         coInvoke(f1, f2);
         result = f1.result + f2.result;
}
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```

#### Cost Model (General)

#### Compositional:

Work: total number of operations

- costs are added across parallel calls

Span: depth/critical path of the computation

- Maximum span is taken across forked calls

Parallelism = Work/Span

- Approximately # of processors that can be effectively used.

#### Combining costs (Nested Parallelism)

#### Combining for parallel for:

$$W_{\text{pexp}}(\text{pfor }...) = \sum_{i=0}^{n-1} W_{\text{exp}}(f(i))$$
 work

$$D_{\text{pexp}}(\text{pfor ...}) = \max_{i=0}^{n-1} D_{\text{exp}}(f(i))$$
 span

## Why Work and Span

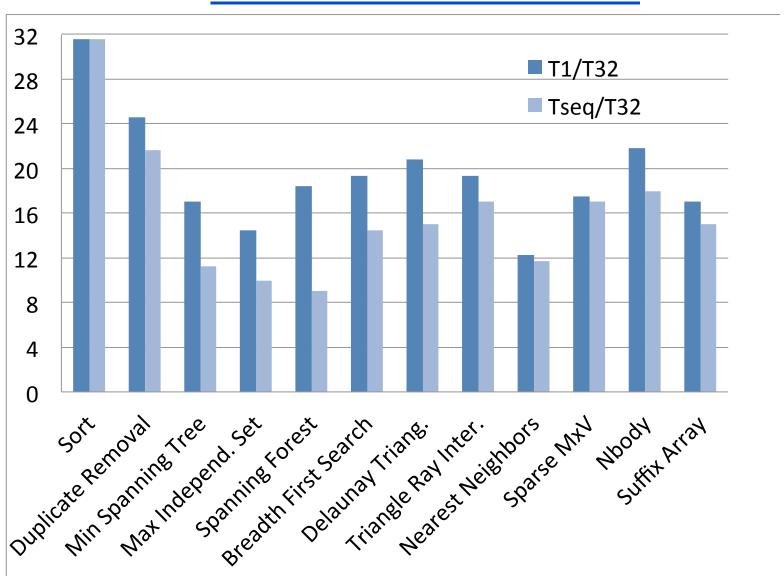
Simple measures that give us a good sense of efficiency (work) and scalability (span).

Can schedule in O(W/P + D) time on P processors.

This is within a constant factor of optimal.

#### Goals in designing an algorithm

- 1. Work should be about the same as the sequential running time. When it matches asymptotically we say it is work efficient.
- 2. Parallelism (W/D) should be polynomial.  $O(n^{1/2})$  is probably good enough



# Styles of Parallel Programming

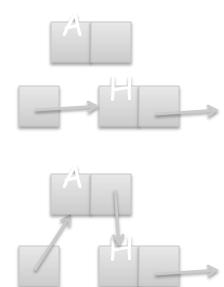
Data parallelism/Bulk Synchronous/SPMD
Nested parallelism: what we covered
Message passing
Futures (other pipelined parallelism)
General Concurrency

# Parallelism vs. Concurrency

- Parallelism: using multiple processors/cores running at the same time. Property of the machine
- Concurrency: non-determinacy due to interleaving threads. Property of the application.

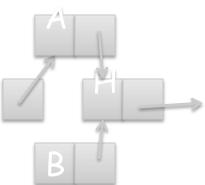
|             |          | Concurrency               |                     |
|-------------|----------|---------------------------|---------------------|
|             |          | sequential                | concurrent          |
| Parallelism | serial   | Traditional programming   | Traditional OS      |
|             | parallel | Deterministic parallelism | General parallelism |

```
struct link {int v; link* next;}
struct stack {
  link* headPtr;
  void push(link* a) {
    a->next = headPtr;
    headPtr = a; }
  link* pop() {
    link* h = headPtr;
    if (headPtr != NULL)
      headPtr = headPtr->next:
    return h;}
```

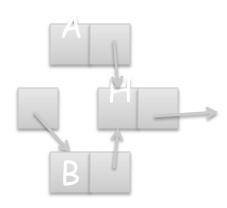


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  void push(link* a) {
    a->next = headPtr;
    headPtr = a; }
  link* pop() {
    link* h = headPtr;
    if (headPtr != NULL)
      headPtr = headPtr->next:
    return h;}
```

```
struct link {int v; link* next;}
struct stack {
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  void push(link* a) {
    a->next = headPtr;
    headPtr = a; }
  link* pop() {
    link* h = headPtr;
    if (headPtr != NULL)
      headPtr = headPtr->next:
    return h;}
```



```
struct link {int v; link* next;}
struct stack {
  link* headPtr;
  void push(link* a) {
    a->next = headPtr;
    headPtr = a; }
  link* pop() {
    link* h = headPtr;
    if (headPtr != NULL)
      headPtr = headPtr->next:
    return h;}
```



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#### <u>CAS</u>

```
bool CAS(ptr* addr, ptr a, ptr b) {
  atomic {
    if (*addr == a) {
        *addr = b;
        return 1;
    } else
        return 0;
}
```

A built in instruction on most processors: CMPXCHG8B - 8 byte version for x86 CMPXCHG16B - 16 byte version

```
struct stack {
  link* headPtr;
  void push(link* a) {
    do {
      link* h = headPtr;
      a->next = h;
    while (!CAS(&headPtr, h, a)); }
  link* pop() {
    do {
      link* h = headPtr;
      if (h == NULL) return NULL;
      link* nxt = h->next;
    while (!CAS(&headPtr, h, nxt))}
    return h;}
```

```
struct stack {
  link* headPtr;
  void push(link* a) {
    do {
      link* h = headPtr;
   a->next = h;
    while (!CAS(&headPtr, h, a)); }
  link* pop() {
    do {
      link* h = headPtr;
      if (h == NULL) return NULL;
      link* nxt = h->next;
    while (!CAS(&headPtr, h, nxt))}
    return h;}
```

```
struct stack {
  link* headPtr;
  void push(link* a) {
    do {
      link* h = headPtr;
    a->next = h;
    while (!CAS(&headPtr, h, a)); }
  link* pop() {
    do {
      link* h = headPtr;
      if (h == NULL) return NULL;
      link* nxt = h->next;
    while (!CAS(&headPtr, h, nxt))}
    return h;}
```

```
struct stack {
  link* headPtr;
  void push(link* a) {
    do {
      link* h = headPtr;
      a->next = h;
   while (!CAS(&headPtr, h, a)); }
  link* pop() {
    do {
      link* h = headPtr;
      if (h == NULL) return NULL;
      link* nxt = h->next;
    while (!CAS(&headPtr, h, nxt))}
    return h;}
```

```
struct stack {
  link* headPtr;
  void push(link* a) {
    do {
      link* h = headPtr;
      a->next = h;
  while (!CAS(&headPtr, h, a)); }
  link* pop() {
    do {
      link* h = headPtr;
      if (h == NULL) return NULL;
      link* nxt = h->next;
    while (!CAS(&headPtr, h, nxt))}
    return h;}
```

```
P1 : x = s.pop(); y = s.pop(); s.push(x);
P2 : z = s.pop();
```

```
Before: B C D
```

After: B C

P2: h = headPtr;

P2: nxt = h->next;

P1: everything

The ABA problem

P2: CAS(&headPtr,h,nxt)

Can be fixed with counter and 2CAS, but...

```
struct link {int v; link* next;}
struct stack {
  link* headPtr;
  void push(link* a) {
    atomic {
      a->next = headPtr;
      headPtr = a; }}
  link* pop() {
    atomic {
      link* h = headPtr;
      if (headPtr != NULL)
        headPtr = headPtr->next;
      return h; } }
```

```
void swapTop(stack s) {
  link* x = s.pop();
  link* y = s.pop();
  push(x);
  push(y);
}
```

Queues are trickier than stacks.

# Styles of Parallel Programming

Data parallelism/Bulk Synchronous/SPMD Nested parallelism: what we covered Message passing

Futures (other pipelined parallelism)

General Concurrency

# Futures: Example

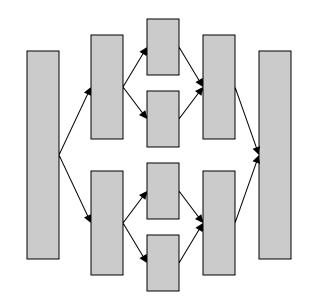
```
fun quickSort S = let
  fun qs([], rest) = rest
    | qs(h::T, rest) =
       let
         val L1 = filter (fn b => b < a) T</pre>
         val L2 = filter (fn b => b >= a) T
       in
         qs(L1,(a::(qs(L2, rest)))
       end
  fun filter(f,[]) = []
      filter(f,h::T) =
        if f(h) then(h::filter(f,T))
        else filter(f,T)
in
 qs(S,[])
15-210
end
```

# Futures: Example

```
fun quickSort S = let
  fun qs([], rest) = rest
    | qs(h::T, rest) =
       let
          val L1 = filter (fn b => b < a) T</pre>
          val L2 = filter (fn b => b >= a) T
        in
          qs(L1, <u>future</u>(a::(qs(L2, rest)))
       end
  fun filter(f,[]) = []
      filter(f,h::T) =
         if f(h) then <u>future</u>(h::filter(f,T))
         else filter(f,T)
in
 qs(S,[])
15-210
end
```

# Quicksort: Nested Parallel

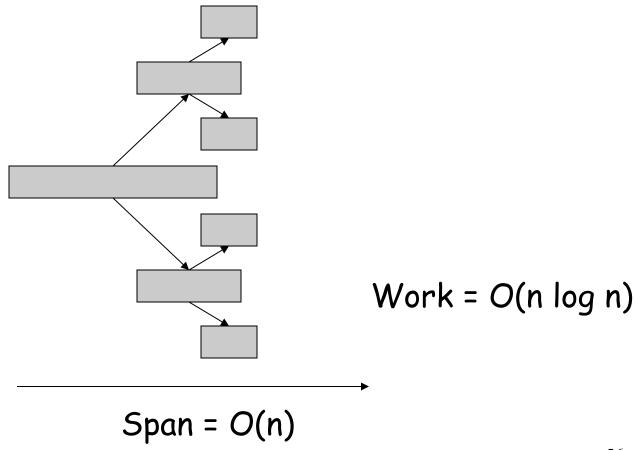
#### Parallel Partition and Append



Work = O(n log n)

Span = 
$$O(lg^2 n)$$

# Quicksort: Futures

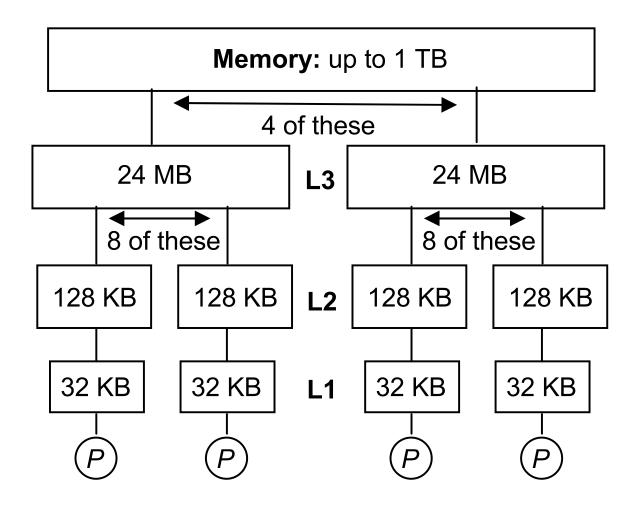


# Styles of Parallel Programming

Data parallelism/Bulk Synchronous/SPMD

- \* Nested parallelism: what we covered Message passing
- \* Futures (other pipelined parallelism)
- \* General Concurrency

#### Xeon 7500



#### Question

How do we get nested parallel programs to work well on a fixed set of processors? Programs say nothing about processors.

Answer: good schedulers

# Greedy Schedules

"Speedup versus Efficiency in Parallel Systems", Eager, Zahorjan and Lazowska, 1989

For any greedy schedule:

Efficiency = 
$$\frac{W}{T_P} \ge \frac{PW}{W + D(P - 1)}$$

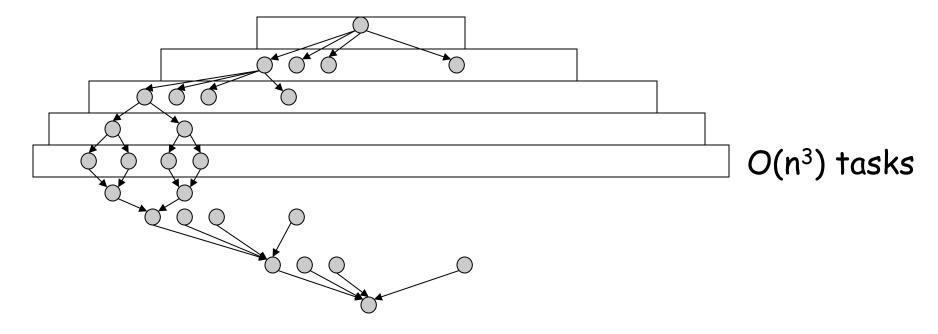
Parallel Time = 
$$T_P \le \frac{W}{P} + D$$

# Types of Schedulers

Bread-First Schedulers
Work-stealing Schedulers
Depth-first Schedulers

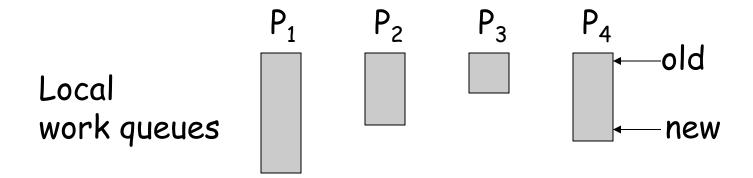
# Breadth First Schedules

Most naïve schedule. Used by most implementations of P-threads.



Bad space usage, bad locality

# Work Stealing



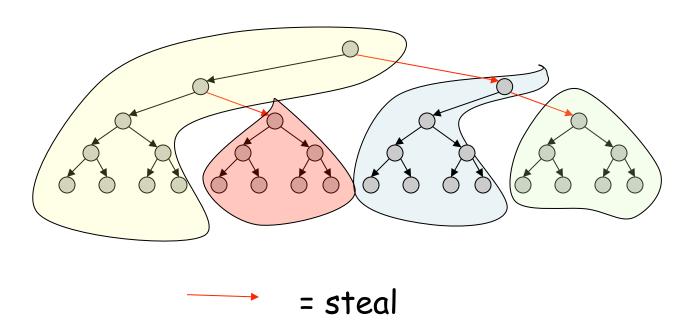
push new jobs on "new" end pop jobs from "new" end

If processor runs out of work, then "steal" from another "old" end

Each processor tends to execute a sequential part of the computation.

# Work Stealing

Tends to schedule "sequential blocks" of tasks



# Work Stealing Theory

```
For strict computations
Blumofe and Leiserson, 1999
# of steals = O(PD)
Space = O(PS_1)  S_1 is the sequential space
Acar, Blelloch and Blumofe, 2003
# of cache misses on distributed caches
M_1 + O(CPD)
M_1 = sequential misses, C = cache size
```

# Work Stealing Practice

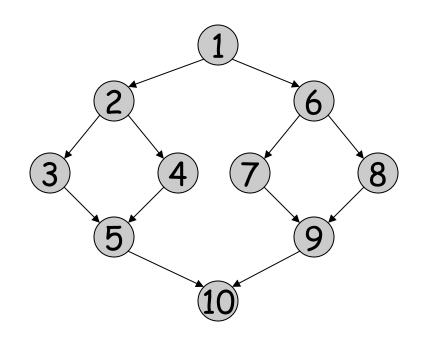
#### Used in Cilk Scheduler

- Small overheads because common case of pushing/popping from local queue can be made fast (with good data structures and compiler help).
- No contention on a global queue
- Has good distributed cache behavior
- Can indeed require O(S<sub>1</sub>P) memory

Used in X10 scheduler, and others

# Parallel Depth First Schedules (P-DFS)

List scheduling based on Depth-First ordering



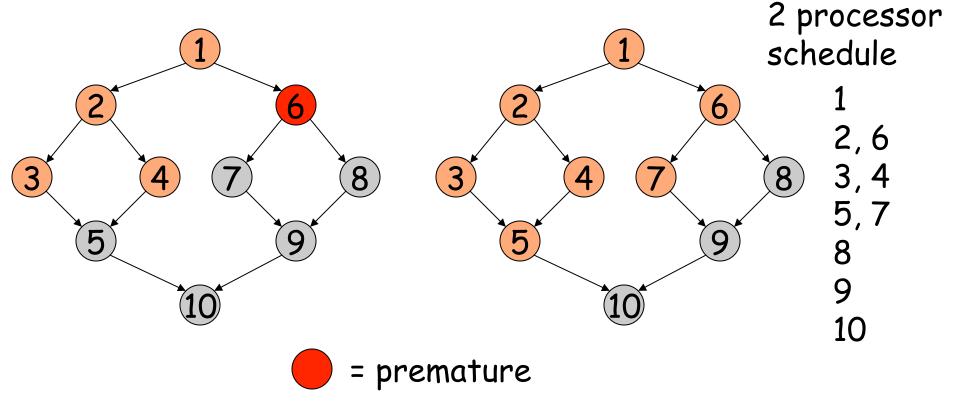
2 processor schedule

1 2,6 3,4 5,7 8 9

For strict computations a shared stack implements a P-DFS

# "Premature task" in P-DFS

A running task is premature if there is an earlier sequential task that is not complete



# P-DFS Theory

Blelloch, Gibbons, Matias, 1999

For any computation:

Premature nodes at any time = O(PD)

Space =  $S_1 + O(PD)$ 

Blelloch and Gibbons, 2004

With a shared cache of size  $C_1$  + O(PD) we have  $M_p$  =  $M_1$ 

# P-DFS Practice

Experimentally uses less memory than work stealing and performs better on a shared cache.

Requires some "coarsening" to reduce overheads

#### Conclusions

- lots of parallel languages
- lots of parallel programming styles
- high-level ideas cross between languages and styles
- scheduling is important