

CSE 332: Data Structures & Parallelism Lecture 15: Analysis of Fork-Join Parallel Programs

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Outline

Done:

- How to use **fork** and **join** to write a parallel algorithm
- Why using divide-and-conquer with lots of small tasks is best
 - Combines results in parallel

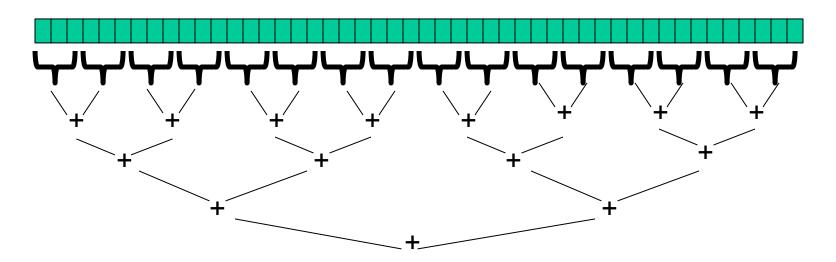
Now:

- More examples of simple parallel programs
- Arrays & balanced trees support parallelism better than linked lists
- Asymptotic analysis for fork-join parallelism
- Amdahl's Law

What else looks like this?

Saw summing an array went from *O*(*n*) sequential to *O*(**log** *n*) parallel (assuming **a lot** of processors and very large *n*)

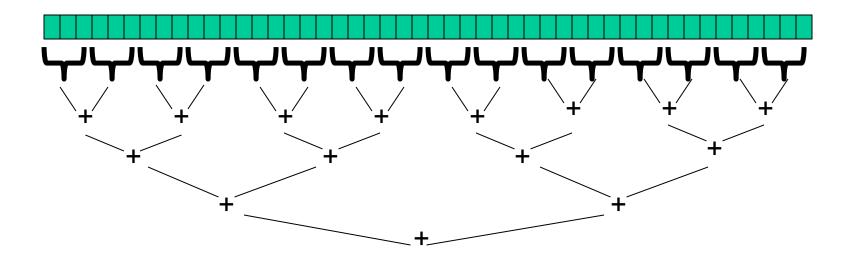
- Exponential speed-up in theory $(n / \log n \text{ grows exponentially})$



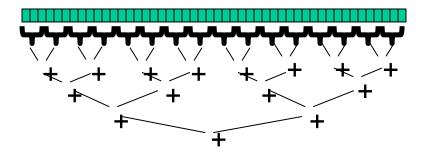
 Anything that can use results from two halves and merge them in O(1) time has the same property...

Extending Parallel Sum

- We can tweak the 'parallel sum' algorithm to do all kinds of things; just specify 2 parts (usually)
 - Describe how to compute the result at the 'cut-off' (Sum: Iterate through sequentially and add them up)
 - Describe how to merge results (Sum: Just add 'left' and 'right' results)



Examples



- Parallelization (for some algorithms)
 - Describe how to compute result at the 'cut-off'
 - Describe how to merge results
- How would we do the following (assuming data is given as an array)?
 - 1. Maximum or minimum element
 - 2. Is there an element satisfying some property (e.g., is there a 17)?
 - 3. Left-most element satisfying some property (e.g., first 17)
 - 4. Smallest rectangle encompassing a number of points
 - 5. Counts; for example, number of strings that start with a vowel
 - 6. Are these elements in sorted order?

Reductions

- This class of computations are called reductions
 - We 'reduce' a large array of data to a single item
 - Produce single answer from collection via an associative operator
 - Examples: max, count, leftmost, rightmost, sum, product, ...
- Note: Recursive results don't have to be single numbers or strings. They can be arrays or objects with multiple fields.
 - Example: create a Histogram of test results from a much larger array of actual test results
- While many can be parallelized due to nice properties like associativity of addition, some things are inherently sequential
 - How we process arr[i] may depend entirely on the result of processing arr[i-1]

Even easier: Maps (Data Parallelism)

- A map operates on each element of a collection independently to create a new collection of the same size
 - No combining results
 - For arrays, this is so trivial some hardware has direct support
- Canonical example: Vector addition

```
int[] vector add(int[] arr1, int[] arr2){
    assert (arr1.length == arr2.length);
    result = new int[arr1.length];
    FORALL(i=0; i < arr1.length; i++) {
        result[i] = arr1[i] + arr2[i];
    }
    return result;
}</pre>
```

Maps in ForkJoin Framework

```
class VecAdd extends RecursiveAction {
  int lo; int hi; int[] res; int[] arr1; int[] arr2;
 VecAdd(int l, int h, int[] r, int[] a1, int[] a2) { ... }
 protected void compute() {
    if(hi - lo < SEQUENTIAL CUTOFF) {</pre>
      for(int i=lo; i < hi; i++)</pre>
        res[i] = arr1[i] + arr2[i];
    } else {
      int mid = (hi+lo)/2;
      VecAdd left = new VecAdd(lo,mid,res,arr1,arr2);
      VecAdd right= new VecAdd(mid,hi,res,arr1,arr2);
      left.fork();
      right.compute();
      left.join();
}
static final ForkJoinPool POOL = new ForkJoinPool();
int[] add(int[] arr1, int[] arr2) {
  assert (arr1.length == arr2.length);
  int[] ans = new int[arr1.length];
  POOL.invoke(new VecAdd(0,arr.length,ans,arr1,arr2);
  return ans;
```

Maps and reductions

Maps and reductions: the "workhorses" of parallel programming

- By far the two most important and common patterns
 - Two more-advanced patterns in next lecture
- Learn to recognize when an algorithm can be written in terms of maps and reductions
- Use maps and reductions to describe (parallel) algorithms
- Programming them becomes "trivial" with a little practice
 - Exactly like sequential for-loops seem second-nature

Map vs reduce in ForkJoin framework

- In our examples:
- Reduce:
 - Parallel-sum extended RecursiveTask
 - Result was returned from compute()
- Map:
 - Class extended was RecursiveAction
 - Nothing returned from compute()
 - In the above code, the 'answer' array was passed in as a parameter
- Doesn't *have* to be this way
 - Map can use RecursiveTask to, say, return an array
 - Reduce could use RecursiveAction; depending on what you're passing back via RecursiveTask, could store it as a class variable and access it via 'left' or 'right' when done

Digression: MapReduce on clusters

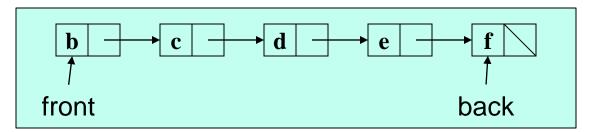
- You may have heard of Google's "map/reduce"
 - Or the open-source version Hadoop
- Idea: Perform maps/reduces on data using many machines
 - The system takes care of distributing the data and managing fault tolerance
 - You just write code to map one element and reduce elements to a combined result
- Separates how to do recursive divide-and-conquer from what computation to perform
 - Old idea in higher-order functional programming transferred to large-scale distributed computing
 - Complementary approach to declarative queries for databases

Trees

- Maps and reductions work just fine on balanced trees
 - Divide-and-conquer each child rather than array sub-ranges
 - Correct for unbalanced trees, but won't get much speed-up
- Example: minimum element in an <u>unsorted</u> but balanced binary tree in O(log n) time given enough processors
- How to do the sequential cut-off?
 - Store number-of-descendants at each node (easy to maintain)
 - Or could approximate it with, e.g., AVL-tree height

Linked lists

- Can you parallelize maps or reduces over linked lists?
 - Example: Increment all elements of a linked list
 - Example: Sum all elements of a linked list
 - Parallelism still beneficial for expensive per-element operations



- Once again, data structures matter!
- For parallelism, balanced trees generally better than lists so that we can get to all the data exponentially faster O(log n) vs. O(n)
 - Trees have the same flexibility as lists compared to arrays (in terms of say inserting an item in the middle of the list)

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Analyzing algorithms

- How to measure efficiency?
 - Want asymptotic bounds
 - Want to analyze the algorithm without regard to a specific number of processors
 - The key "magic" of the ForkJoin Framework is getting expected run-time performance asymptotically optimal for the available number of processors
 - So we can analyze algorithms assuming this guarantee

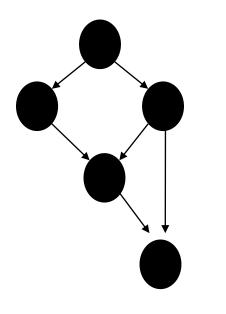
Work and Span

Let T_P be the running time if there are P processors available Two key measures of run-time:

- Work: How long it would take 1 processor = T₁
 - Just "sequentialize" the recursive forking
 - Cumulative work that all processors must complete
- Span: How long it would take infinity processors = T_{∞}
 - The hypothetical ideal for parallelization
 - This is the longest "dependence chain" in the computation
 - Example: O(log n) for summing an array
 - Notice in this example having > n/2 processors is no additional help
 - Also called "critical path length" or "computational depth"

The DAG (Directed Acyclic Graph)

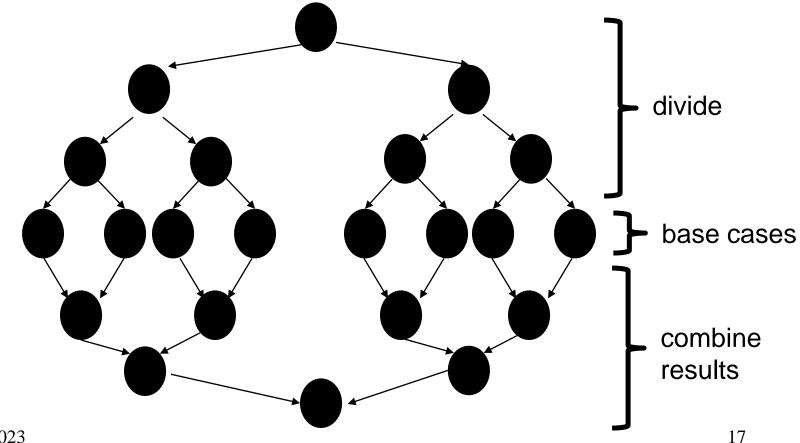
- A program execution using **fork** and **join** can be seen as a DAG
- [A DAG is a graph that is <u>directed</u> (edges have direction (arrows)), and those arrows do not create a <u>cycle</u> (ability to trace a path that starts and ends at the same node).]
 - Nodes: Pieces of work
 - Edges: Source must finish before destination starts



- A fork "ends a node" and makes two outgoing edges
 - New thread
 - Continuation of current thread
- A join "ends a node" and makes a node with two incoming edges
 - Node just ended
 - Last node of thread joined on

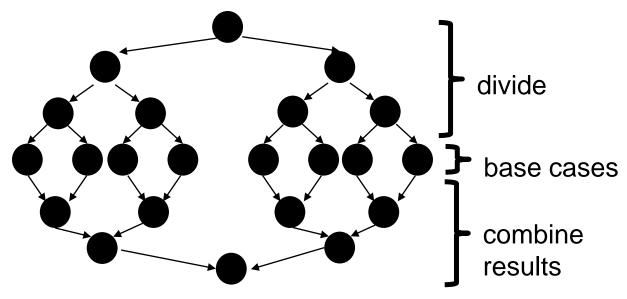
Our simple examples

- fork and join are very flexible, but divide-and-conquer maps and reductions use them in a very basic way:
 - A tree on top of an upside-down tree



Our simple examples, in more detail

Our fork and join frequently look like this:



In this context, the span (T_{∞}) is:

- •The longest dependence-chain; longest 'branch' in parallel 'tree'
- •Example: O(log n) for summing an array; we halve the data down to our
- cut-off, then add back together; O(log n) steps, O(1) time for each
- •Also called "critical path length" or "computational depth"

More interesting DAGs?

- The DAGs are not always this simple
- Example:
 - Suppose combining two results might be expensive enough that we want to parallelize each one
 - Then each node in the inverted tree on the previous slide would itself expand into another set of nodes for that parallel computation

Connecting to performance

- Recall: T_P = running time if there are P processors available
- Work = T_1 = sum of run-time of all nodes in the DAG
 - That lonely processor does everything
 - Any topological sort is a legal execution
 - O(n) for simple maps and reductions
- Span = T_{∞} = sum of run-time of all nodes on the most-expensive path in the DAG
 - Note: costs are on the nodes not the edges
 - Our infinite army can do everything that is ready to be done, but still has to wait for earlier results
 - O(log n) for simple maps and reductions

Definitions

A couple more terms:

- Speed-up on P processors: T₁ / T_P
- If speed-up is **P** as we vary **P**, we call it perfect linear speed-up
 - Perfect linear speed-up means doubling **P** halves running time
 - Usually our goal; hard to get in practice
- Parallelism is the maximum possible speed-up: T_1 / T_{∞}
 - At some point, adding processors won't help
 - What that point is depends on the span

Parallel algorithms is about decreasing span without increasing work too much

Optimal T_P: Thanks ForkJoin library!

- So we know T_1 and T_{∞} but we want T_P (e.g., P=4)
- Ignoring memory-hierarchy issues (caching), **T**_P can't beat
 - **T**₁ / **P** why not?
 - $-\mathbf{T}_{\infty}$ why not?
- So an *asymptotically* optimal execution would be:

 $T_{P} = O((T_{1} / P) + T_{\infty})$

- First term dominates for small P, second for large P
- The ForkJoin Framework gives an *expected-time guarantee* of asymptotically optimal!
 - Expected time because it flips coins when *scheduling*
 - How? For an advanced course (few need to know)
- Guarantee requires a few assumptions about your code...
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Division of responsibility

- Our job as ForkJoin Framework users:
 - Pick a good algorithm, write a program
 - When run, program creates a DAG of things to do
 - Make all the nodes a small-ish and approximately equal amount of work
- The framework-writer's job:
 - Assign work to available processors to avoid idling
 - Let framework-user ignore all scheduling issues
 - Keep constant factors low
 - Give the expected-time optimal guarantee assuming framework-user did his/her job

$$T_{P} = O((T_{1} / P) + T_{\infty})$$

Examples

$T_{P} = O((T_{1} / P) + T_{\infty})$

- In the algorithms seen so far (e.g., sum an array):
 - $\mathbf{T}_1 = O(n)$
 - $\mathbf{T}_{\infty} = O(\log n)$
 - So expect (ignoring overheads): $T_P = O(n/P + \log n)$
- Suppose instead:
 - $\mathbf{T}_1 = O(n^2)$
 - $\mathbf{T}_{\infty} = O(n)$
 - So expect (ignoring overheads): $T_P = O(n^2/P + n)$

And now for the bad news...

- So far: talked about a parallel program in terms of work and span
- In practice, it's common that your program has:

a) parts that **parallelize well**:

- Such as maps/reduces over arrays and trees
- b) ...and parts that **don't parallelize at all:**
- Such as reading a linked list, getting input, or just doing computations where each step needs the results of previous step
- These unparallelized parts can turn out to be a big bottleneck, which brings us to Amdahl's Law ...

Amdahl's Law (mostly bad news)

Let the *work* (time to run on 1 processor) be 1 unit time

Let S be the portion of the execution that can't be parallelized

Suppose we get perfect linear speedup on the parallel portion

Then: **T**_P = **S** + (1-**S**)/**P**

So the overall speedup with P processors is (Amdahl's Law): $T_1 / T_P = 1 / (S + (1-S)/P)$

And the parallelism (infinite processors) is:

 $T_1 / T_{\infty} = 1 / S$

Amdahl's Law Example

Suppose: $T_1 = S + (1-S) = 1$ (aka total program execution time) $T_1 = 1/3 + 2/3 = 1$ $T_1 = 33 \text{ sec} + 67 \text{ sec} = 100 \text{ sec}$

Time on P processors: $T_P = S + (1-S)/P$

So:
$$T_P = 33 \text{ sec} + (67 \text{ sec})/P$$

 $T_3 = 33 \text{ sec} + (67 \text{ sec})/3 =$

Why such bad news?

 $T_1 / T_P = 1 / (S + (1-S)/P)$ $T_1 / T_{\infty} = 1 / S$

- Suppose 33% of a program is sequential
 - Then a billion processors won't give a speedup over 3!!!
- No matter how many processors you use, your speedup is bounded by the sequential portion of the program.

The future and Amdahl's Law

 Speedup:
 $T_1 / T_P = 1 / (S + (1-S)/P)$

 Max Parallelism:
 $T_1 / T_\infty = 1 / S$

- Suppose you miss the good old days (1980-2005) where 12ish years was long enough to get 100x speedup
 - Now suppose in 12 years, clock speed is the same but you get 256 processors instead of 1
 - What portion of the program must be parallelizable to get 100x speedup?

All is not lost

Amdahl's Law is a bummer!

- Unparallelized parts become a bottleneck very quickly
- But it doesn't mean additional processors are worthless
- We can find new parallel algorithms
 - Some things that seem entirely sequential turn out to be parallelizable
 - Eg. How can we parallelize the following?
 - Take an array of numbers, return the 'running sum' array:

input	6	4	16	10	16	14	2	8
output	6	10	26	36	52	66	68	76

- At a glance, not sure; we'll explore this shortly
- We can also change the problem we're solving or do new things
 - Example: Video games use tons of parallel processors
 - They are not rendering 10-year-old graphics faster
 - They are rendering richer environments and more beautiful (terrible?) monsters

Moore and Amdahl



- Moore's "Law" is an observation about the progress of the semiconductor industry
 - Transistor density doubles roughly every 18 months
- Amdahl's Law is a mathematical theorem
 - Diminishing returns of adding more processors
- Both are incredibly important in designing computer systems