CSE 332: Data Abstractions

Lecture 18: 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
• Some Java and ForkJoin Framework specifics
  – More pragmatics (e.g., installation) in separate notes

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 ($n / \log n$ 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 (proj3)
  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

```java
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;
}
```
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) {
            for (int i = lo; i < hi; i++)
                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 fjPool = new ForkJoinPool();
int[] add(int[] arr1, int[] arr2) {
    assert (arr1.length == arr2.length);
    int[] ans = new int[arr1.length];
    fjPool.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 *unsorted* 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)
Analyzing algorithms

• Like all algorithms, parallel algorithms should be:
  – Correct
  – Efficient

• For our algorithms so far, correctness is “obvious” so we’ll focus on 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 = $T_1$
  - Just “sequentialize” the recursive forking

- **Span**: How long it would take infinity processors = $T_\infty$
  - 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

• A program execution using fork and join can be seen as a DAG
  – **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

![Diagram of a tree structure with arrows indicating divide, base cases, and combine results.](image-url)
Our simple examples, in more detail

Our fork and join frequently look like this:

In this context, the span ($T_\infty$) 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: $T_P = \text{running time if there are } P \text{ processors available}$

• Work $= T_1 = \text{sum of run-time of all nodes in the DAG}$
  – That lonely processor does everything
  – Any topological sort is a legal execution
  – $O(n)$ for simple maps and reductions

• Span $= T_\infty = \text{sum of run-time of all nodes on the most-expensive path in the DAG}$
  – 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: $T_1 / T_P$

• 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: $T_1 / T_\infty$
  – 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 $T_\infty$ but we want $T_P$ (e.g., $P=4$)
- Ignoring memory-hierarchy issues (caching), $T_P$ can’t beat
  - $T_1 / P$ why not?
  - $T_\infty$ why not?
- So an asymptotically optimal execution would be:

  $$T_P = O\left(\frac{T_1}{P} + T_\infty\right)$$
  - 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…
**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

\[ T_P = O\left(\frac{T_1}{P} + T_\infty\right) \]
Examples

\[ T_P = O((T_1 / P) + T_\infty) \]

- In the algorithms seen so far (e.g., sum an array):
  - \( T_1 = O(n) \)
  - \( T_\infty = O(\log n) \)
  - So expect (ignoring overheads): \( T_P = O(n/P + \log n) \)

- Suppose instead:
  - \( T_1 = O(n^2) \)
  - \( T_\infty = O(n) \)
  - So expect (ignoring overheads): \( T_P = O(n^2/P + n) \)
Amdahl’s Law (mostly 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
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.

Then:

$$T_1 = S + (1-S) = 1$$

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:

$$T_1 / T_\infty = 1 / S$$
Why such bad news?

\[
\frac{T_1}{T_P} = \frac{1}{S + (1-S)/P} \quad \quad \quad \frac{T_1}{T_\infty} = \frac{1}{S}
\]

- 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.
Amdahl’s Law Example

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

Time on P processors: \[ T_P = S + \frac{(1-S)}{P} \]

So: \[ T_P = 33 \text{ sec} + \frac{(67 \text{ sec})}{P} \]
\[ T_3 = 33 \text{ sec} + \frac{(67 \text{ sec})}{3} = \]
The future and Amdahl’s Law

Speedup: \[
\frac{T_1}{T_P} = \frac{1}{S + \frac{(1-S)}{P}}
\]
Max Parallelism: \[
\frac{T_1}{T_\infty} = \frac{1}{S}
\]

• 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?
The future and Amdahl’s Law

Speedup: \( \frac{T_1}{T_P} = \frac{1}{S + \frac{1-S}{P}} \)

Max Parallelism: \( \frac{T_1}{T_\infty} = \frac{1}{S} \)

• 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?

For 256 processors to get at least 100x speedup, we need

\[ 100 \leq \frac{1}{S + \frac{1-S}{256}} \]

Which means \( S \leq .0061 \) (i.e., 99.4% must be parallelizable)
Plots you have to see

1. Assume 256 processors
   - x-axis: sequential portion $S$, ranging from .01 to .25
   - y-axis: speedup $T_1 / T_P$ (will go down as $S$ increases)

2. Assume $S = .01$ or .1 or .25 (three separate lines)
   - x-axis: number of processors $P$, ranging from 2 to 32
   - y-axis: speedup $T_1 / T_P$ (will go up as $P$ increases)

Do this as a homework problem!
   - Chance to use a spreadsheet or other graphing program
   - Compare against your intuition
   - A picture is worth 1000 words, especially if you made it
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