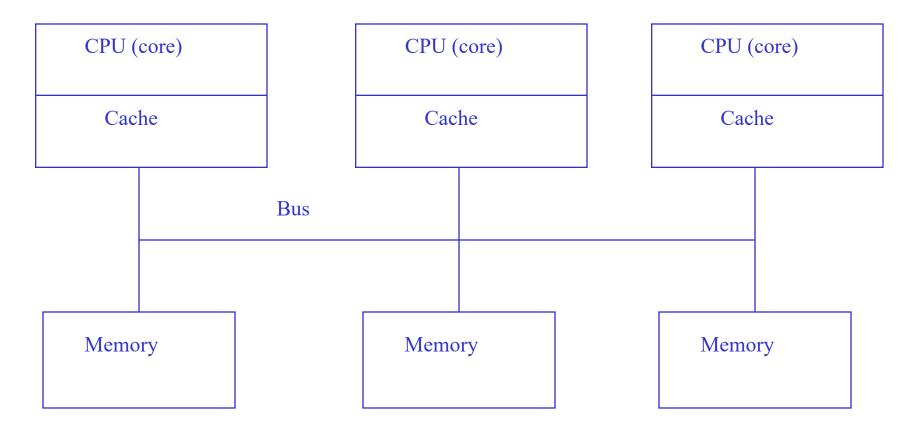
### Parallel Scientific Programming

#### • Definitions

- Speedup:  $T_s/T_p$ , where  $T_s$  is the sequential time and  $T_p$  is the time when using *p* cores
  - "Perfect speedup" is *p*, which really should be called "linear speedup"
  - Typically, speedup is less than *p*---but it can be larger because of memory hierarchy effects
- Efficiency: Speedup/p
  - Intuition: how well am I using my *p* cores

#### "Superlinear" Speedup?



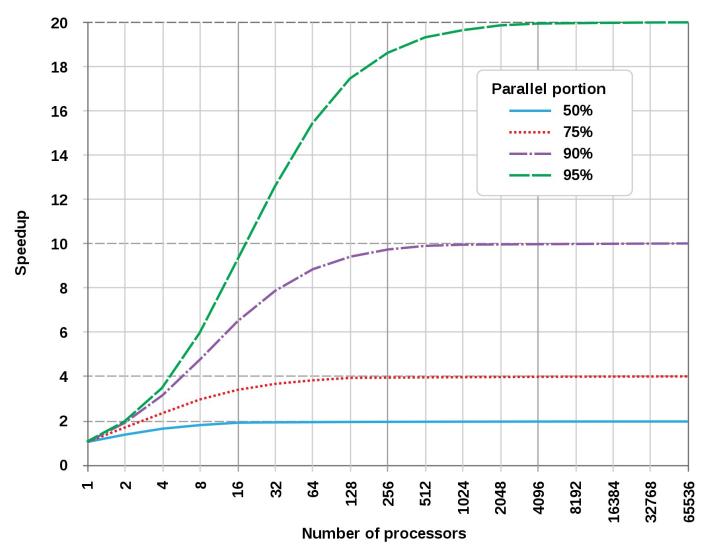
Strictly more cache when more cores are used Can result in fewer cache misses when using more cores Same argument can hold for any level of the memory hierarchy

### Graphs of Speedup and Efficiency (shown on board)

### Parallel Scientific Programming

- Definitions, continued
  - Amdahl's law: if  $T_s$  is sequential time, then:
    - $T_p \approx T_s * (1-f) + (T_s / p) * f$ , where *f* is the fraction of the program that is parallelizable, and *p* is the number of cores.
    - Intuition: the non-parallelizable portion doesn't speed up at all, and the parallelizable part scales linearly
    - Of course, this isn't true in general; one might, for example, have load imbalance in a parallelizable portion---also ignored is process/thread creation, communication, synchronization

#### Amdahl's Law



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#### Example use of Amdahl's Law

- Suppose program takes 50 seconds sequentially, but of that 50 seconds:
  - 5 seconds is initialization that can't be parallelized
  - 5 seconds is finalization that also can't be parallelized
- Then the **maximum speedup** possible is 5!
  - Even if we have an arbitrarily large number of cores!
  - Lesson: try to avoid sequential portions of code

# Different parallel programming styles (end of Chapter 3)

- Iterative: SPMD (e.g., Jacobi iteration)
- Recursive: adaptive quadrature
- Task parallel: independent portions of programs
- Bag of tasks: implementation of recursive, generally, but also can be used for iterative

#### Data Parallel Algorithms

- Execute identical code on different parts of a data structure
  - Usually we mean "SPMD algorithms", which stands for Single Program Multiple Data
  - Data Parallel implies barrier after every instruction (arose from programming SIMD architectures; recall SIMD is Single Instruction Multiple Data)
  - SPMD allows barriers at arbitrary points (arose from MIMD architectures)
    - It is a relaxation of SIMD

## Picture: finding the sum of an array in parallel (SPMD program)

```
Finding the sum of an array in parallel
               (SPMD program)
int sum[n], old[n], a[n] // Array a initialized to arbitrary values
coi := 0 to n-1
 int d = 1
 sum[i] = a[i]
 while (d < n) {
   old[i] = sum[i]
   barrier ←
                                      Why?
   if (i - d \ge 0)
     sum[i] = old[i-d] + sum[i]
   barrier ←
                                      Why?
   d = d * 2
  }
OC
```

#### Picture: Jacobi Iteration