



Parallel software

The burden is on software

- Hardware and compilers can keep up the pace needed.
- From now on...
 - In shared memory programs:
 - Start a single process and fork threads.
 - Threads carry out tasks.
 - In distributed memory programs:
 - Start multiple processes.
 - Processes carry out tasks.

SPMD – single program multiple data

- A SPMD programs consists of a single executable that can behave as if it were multiple different programs through the use of conditional branches.

```
if (I'm thread process i)
    do this;
else
    do that;
```



Writing Parallel Programs

1. Divide the work among the processes/threads
 - (a) so each process/thread gets roughly the **same amount** of work
 - (b) and communication is **minimized**.
2. Arrange for the processes/threads to **synchronize**.
3. Arrange for **communication** among processes/threads.

```
double x[n], y[n];  
...  
for (i = 0; i < n; i++)  
    x[i] += y[i];
```

Shared Memory

- **Dynamic threads**

- Master thread **waits** for work, **forks** new threads, and when threads are done, they **terminate**
- Efficient use of resources, but thread **creation** and **termination** is **time consuming**.

- **Static threads**

- Pool of threads are created and allocated work, but do **not terminate until cleanup**.
- Better performance, but **potential waste** of system resources.

Nondeterminism

...

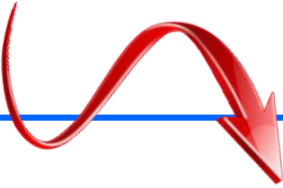
```
printf ( "Thread %d > my_val = %d\n" ,  
        my_rank , my_x ) ;
```

...



Thread 1 > my_val = 19

Thread 0 > my_val = 7



Thread 0 > my_val = 7

Thread 1 > my_val = 19

Nondeterminism

```
my_val = Compute_val ( my_rank ) ;  
x += my_val ;
```

Time	Core 0	Core 1
0	Finish assignment to my_val	In call to Compute_val
1	Load x = 0 into register	Finish assignment to my_val
2	Load my_val = 7 into register	Load x = 0 into register
3	Add my_val = 7 to x	Load my_val = 19 into register
4	Store x = 7	Add my_val to x
5	Start other work	Store x = 19

Nondeterminism

- Race condition
- Critical section
- Mutually exclusive
- Mutual exclusion lock (mutex, or simply lock)

```
my_val = Compute_val ( my_rank ) ;  
Lock(&add_my_val_lock ) ;  
x += my_val ;  
Unlock(&add_my_val_lock ) ;
```


busy-waiting

```
my_val = Compute_val ( my_rank ) ;  
if ( my_rank == 1)  
    while ( ! ok_for_1 ) ; /* Busy-wait loop */  
x += my_val ; /* Critical section */  
if ( my_rank == 0)  
    ok_for_1 = true ; /* Let thread 1 update x */
```

message-passing

```
char message [ 1 0 0 ] ;
```

```
...
```

```
my_rank = Get_rank ( ) ;
```

```
if ( my_rank == 1 ) {
```

```
    printf ( message , "Greetings from process 1" ) ;
```

```
    Send ( message , MSG_CHAR , 100 , 0 ) ;
```

```
} elseif ( my_rank == 0 ) {
```

```
    Receive ( message , MSG_CHAR , 100 , 1 ) ;
```

```
    printf ( "Process 0 > Received: %s\n" , message ) ;
```

```
}
```

Partitioned Global Address Space Languages

```
shared int n = ... ;  
shared double x [ n ] , y [ n ] ;  
private int i , my_first_element , my_last_element ;  
my_first_element = ... ;  
my_last_element = ... ;  
/* Initialize x and y */  
...  
for ( i = my_first_element ; i <= my_last_element ; i++)  
    x [ i ] += y [ i ] ;
```

Input and Output

- In distributed memory programs, only **process 0** will access *stdin*.
- In shared memory programs, only the **master thread or thread 0** will access *stdin*.
- In both distributed memory and shared memory programs all the processes/threads can access *stdout* and *stderr*.

Input and Output

- In most cases only **a single process/thread** will be used for all output to ***stdout*** other than debugging output.
 - because of the indeterminacy of the order of output to ***stdout***
- **Debug output** should always include **the rank or id** of the process/thread that's generating the output.

Input and Output

- Only **a single process/thread** will attempt to access **any single file** other than stdin, stdout, or stderr.
 - E.g., each process/thread can open its own, private file for reading or writing,
- But **no two processes/threads** will open the **same file**.

Performance



Speedup



- Number of cores = p
- Serial run-time = T_{serial}
- Parallel run-time = T_{parallel}

linear speedup

$$T_{\text{parallel}} = T_{\text{serial}} / p$$

Speedup of a parallel program

$$S = \frac{T_{\text{serial}}}{T_{\text{parallel}}}$$

Efficiency of a parallel program

$$E = \frac{S}{p} = \frac{\left(\frac{T_{\text{serial}}}{T_{\text{parallel}}} \right)}{p} = \frac{T_{\text{serial}}}{p \cdot T_{\text{parallel}}}$$

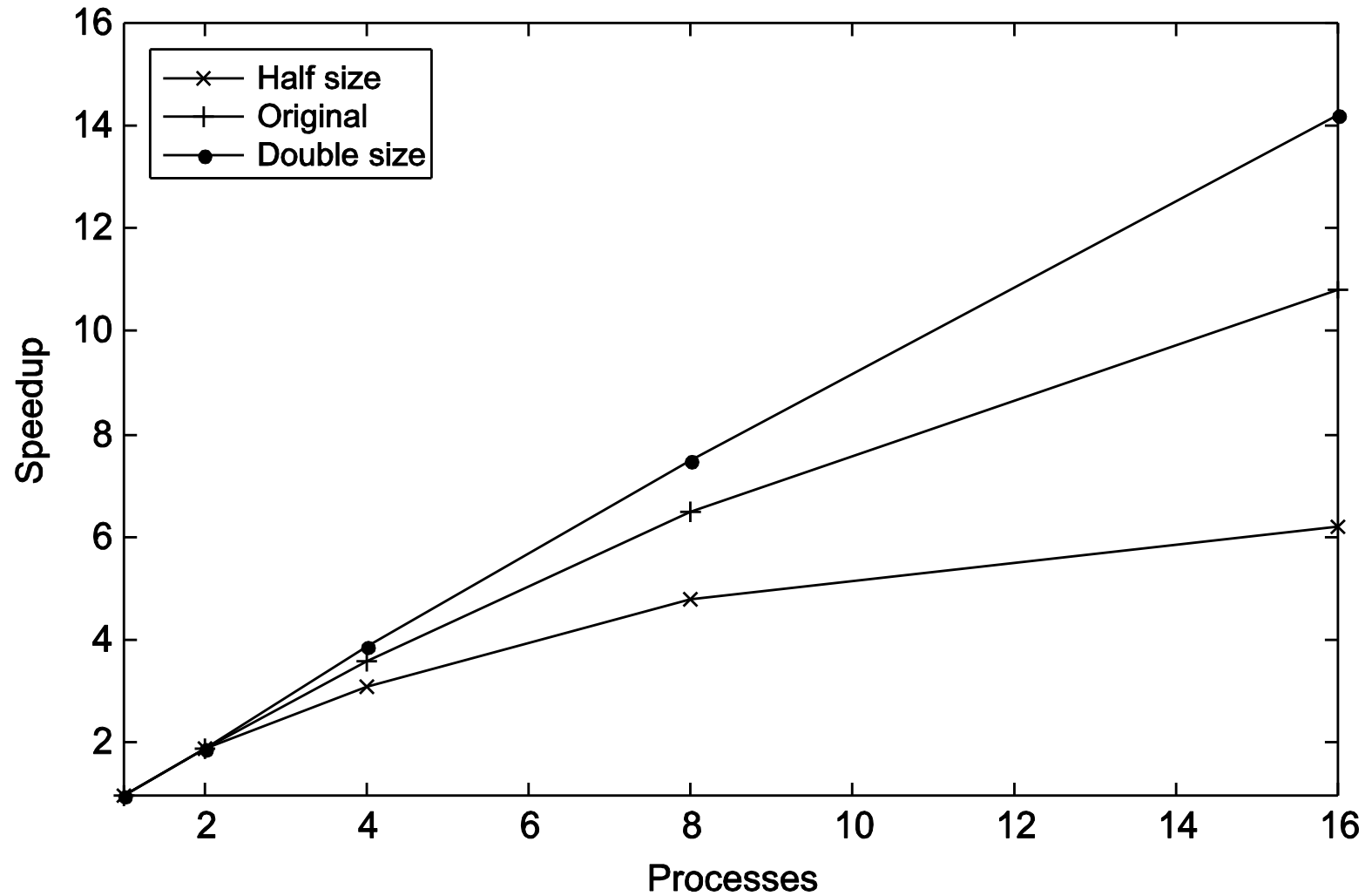
Speedups and efficiencies of a parallel program

p	1	2	4	8	16
S	1.0	1.9	3.6	6.5	10.8
$E = S/p$	1.0	0.95	0.90	0.81	0.68

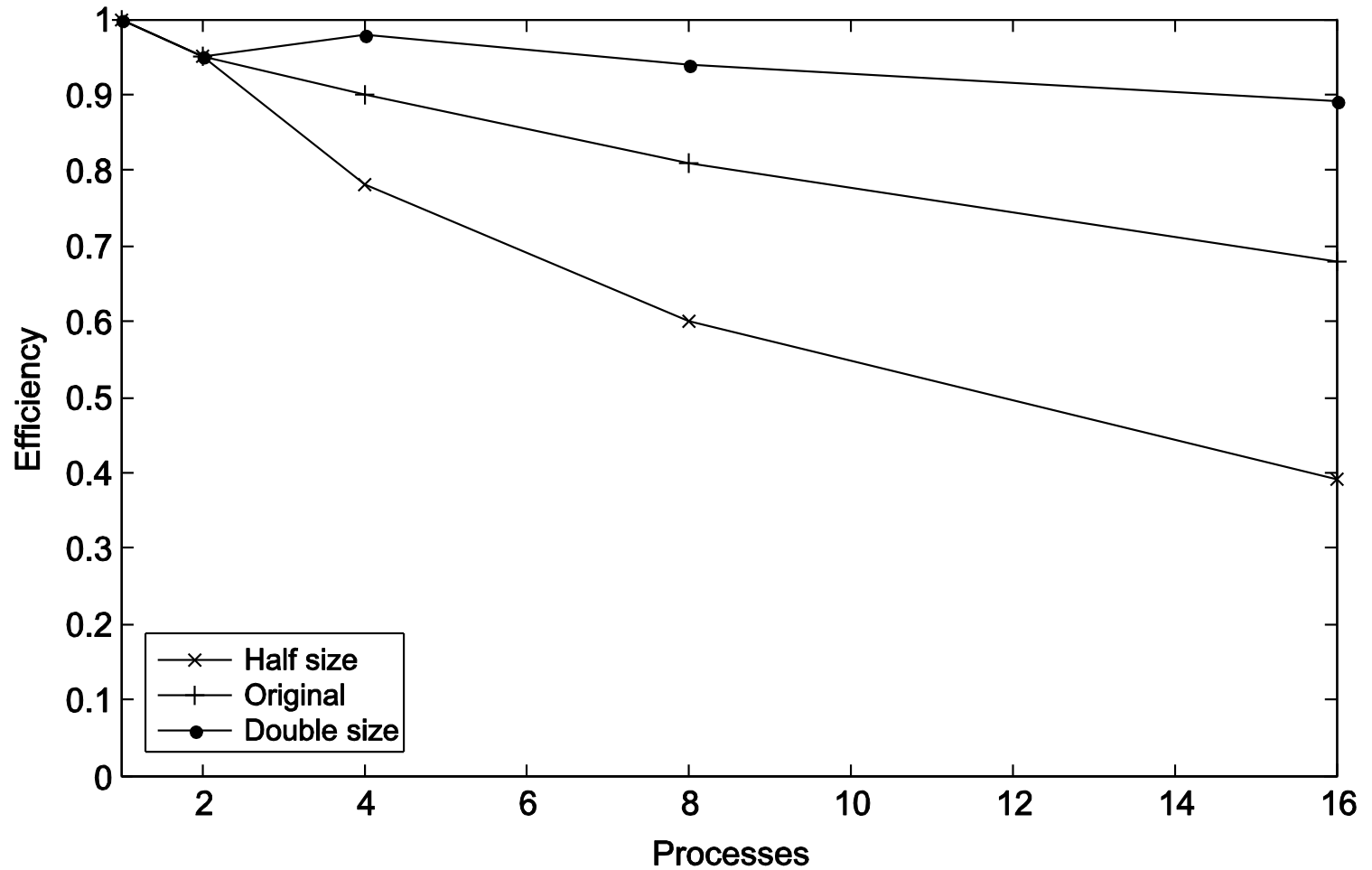
Speedups and efficiencies of parallel program on different problem sizes

	p	1	2	4	8	16
Half	S	1.0	1.9	3.1	4.8	6.2
	E	1.0	0.95	0.78	0.60	0.39
Original	S	1.0	1.9	3.6	6.5	10.8
	E	1.0	0.95	0.90	0.81	0.68
Double	S	1.0	1.9	3.9	7.5	14.2
	E	1.0	0.95	0.98	0.94	0.89

Speedup



Efficiency



Effect of overhead

$$T_{\text{parallel}} = T_{\text{serial}} / p + T_{\text{overhead}}$$

Amdahl's Law

- Unless virtually all of a serial program is parallelized, the possible speedup is going to be very limited — regardless of the number of cores available.



Example

- We can parallelize **90%** of a serial program.
- Parallelization is “perfect” regardless of the number of cores p we use.
- $T_{\text{serial}} = 20$ seconds
- Runtime of parallelizable part is

$$0.9 \times T_{\text{serial}} / p = 18 / p$$

Example (cont.)

- Runtime of “unparallelizable” part is

$$0.1 \times T_{serial} = 2$$

- Overall parallel run-time is

$$T_{parallel} = 0.9 \times T_{serial} / p + 0.1 \times T_{serial} = 18 / p + 2$$

Example (cont.)

- Speed up

$$S = \frac{T_{serial}}{0.9 \times T_{serial} / p + 0.1 \times T_{serial}} = \frac{20}{18 / p + 2}$$

Scalability

- In general, a problem is *scalable* if it can handle ever increasing problem sizes.
- If we increase the number of processes/threads and keep the efficiency fixed *without increasing problem size*, the problem is *strongly scalable*.
- If we keep the efficiency fixed by *increasing the problem size* at the same rate as we increase the number of processes/threads, the problem is *weakly scalable*.

强可扩展性 (Strongly Scalable)

- **定义：**在**问题规模（即任务大小）固定**的情况下，通过增加计算资源（如处理器数量）来**缩短计算时间**。
- **目标：**提高计算速度，使任务更快完成。
- **关键点：****问题规模不变，资源增加，计算时间减少。**
- **理想情况：**资源增加一倍，计算时间减半（线性加速）。
- **例子：**假设你需要对一个固定大小的矩阵（例如 1000×1000 ）进行矩阵乘法运算：
 - 使用 **1 个处理器**，计算时间为 100 秒。
 - 使用 **2 个处理器**，计算时间减少到 50 秒。
 - 使用 **4 个处理器**，计算时间减少到 25 秒。

弱可扩展性 (Weakly Scalable)

- **定义：**在**增加计算资源的同时，问题规模也按比例增加**，目标是保持每个处理器的负载不变，**计算时间基本稳定**。
- **目标：**在资源增加的情况下，能够处理更大规模的问题，同时保持效率。
- **关键点：****问题规模与资源同步增加，计算时间不变**。
- **理想情况：**资源增加一倍，问题规模也增加一倍，计算时间保持不变
- **例子：**假设你需要对一个矩阵进行乘法运算（**对于方阵 $n \times n$ ，运算量为 n^3 次乘法**），但矩阵的大小会随着处理器数量的增加而增加：
 - 使用 **1 个处理器**，计算一个 100×100 的矩阵，时间为 10 秒。
 - 使用 **8 个处理器**，计算一个 200×200 的矩阵，时间仍为 10 秒。
 - 使用**64个处理器**，计算一个 400×400 的矩阵，时间仍为 10 秒。

区别总结

特性	强可扩展性	弱可扩展性
问题规模	固定	随资源增加而增加
资源增加	资源增加，问题规模不变	资源增加，问题规模同步增加
目标	缩短计算时间	处理更大规模的问题，保持计算时间稳定
适用场景	计算密集型任务（如固定规模的科学计算）	数据密集型任务（如大数据处理）

- **强可扩展性**关注的是在固定问题规模下，通过增加资源来加速计算。
- **弱可扩展性**关注的是在资源增加的同时，能够处理更大规模的问题，同时保持计算效率。

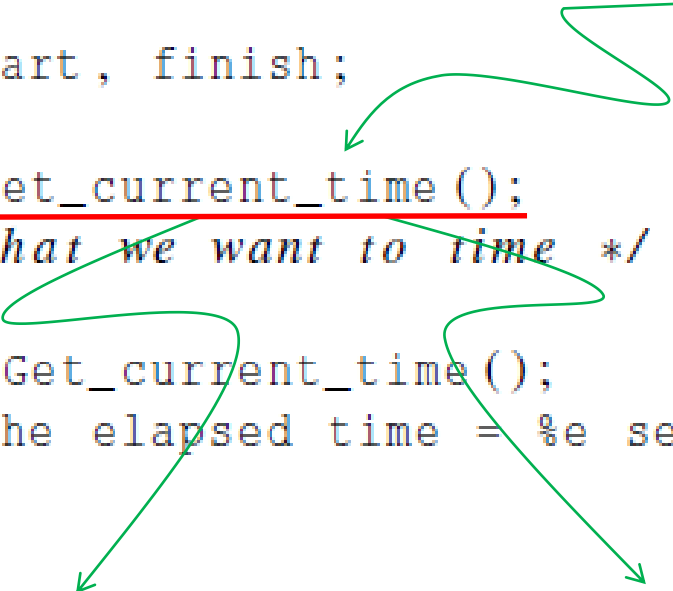
Taking Timings

- What is time?
- Start to finish?
- A program segment of interest?
- CPU time?
- Wall clock time?



Taking Timings

```
double start, finish;  
...  
start = Get_current_time();  
/* Code that we want to time */  
...  
finish = Get_current_time();  
printf("The elapsed time = %e seconds\n", finish-start);
```



theoretical
function

MPI_Wtime

omp_get_wtime

Taking Timings

```
private double start, finish;  
...  
start = Get_current_time();  
/* Code that we want to time */  
...  
finish = Get_current_time();  
printf("The elapsed time = %e seconds\n", finish-start);
```

Taking Timings

```
shared double global_elapsed;  
private double my_start, my_finish, my_elapsed;  
.  
.  
.  
/* Synchronize all processes/threads */  
Barrier();  
my_start = Get_current_time();  
  
/* Code that we want to time */  
.  
.  
.  
  
my_finish = Get_current_time();  
my_elapsed = my_finish - my_start;  
  
/* Find the max across all processes/threads */  
global_elapsed = Global_max(my_elapsed);  
if (my_rank == 0)  
    printf("The elapsed time = %e seconds\n", global_elapsed);
```



Parallel program design

Foster's methodology

1. **Partitioning**: divide the computation to be performed and the data operated on by the computation into small tasks.

The focus here should be on **identifying tasks** that can be **executed in parallel**.

Foster's methodology

2. **Communication**: determine what communication needs to be carried out among the tasks identified in the previous step.



Foster's methodology

3. **Agglomeration or aggregation**: combine tasks and communications identified in the first step into larger tasks.

For example, if task A must be executed before task B can be executed, it may make sense to aggregate them into a single **composite task**.

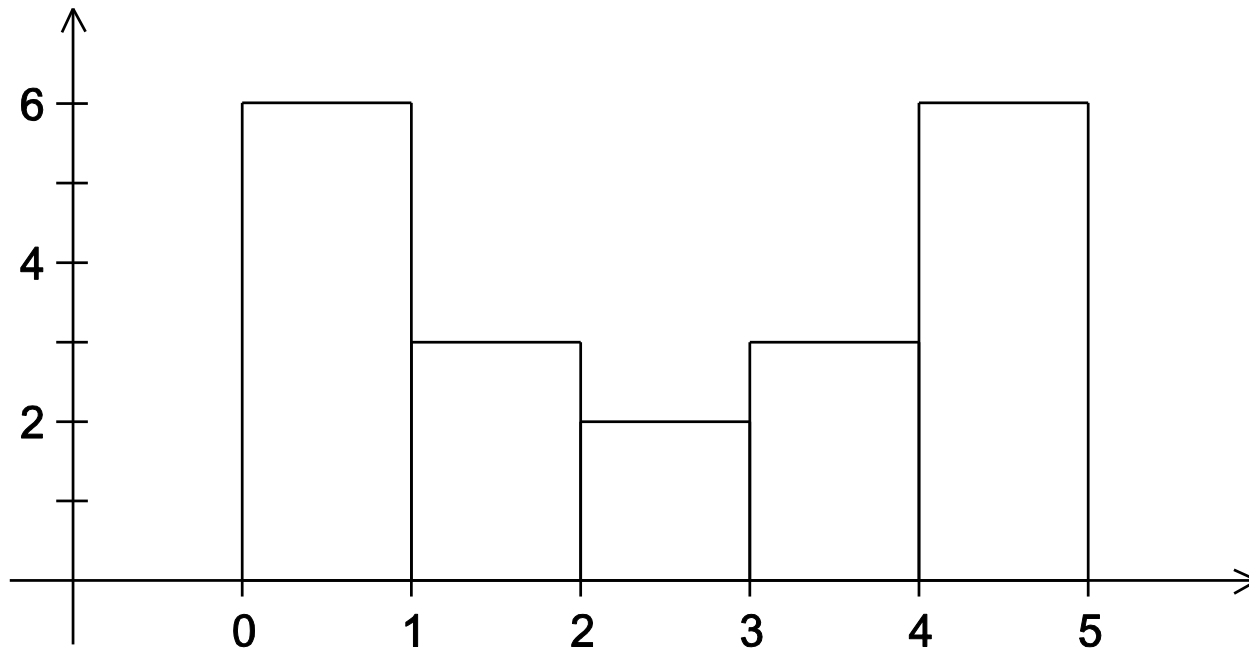
Foster's methodology

4. **Mapping**: assign the composite tasks identified in the previous step to processes/threads.

This should be done so that communication is minimized, and each process/thread gets roughly the same amount of work.

Example - histogram

- 1.3, 2.9, 0.4, 0.3, 1.3, 4.4, 1.7, 0.4, 3.2, 0.3, 4.9, 2.4, 3.1, 4.4, 3.9, 0.4, 4.2, 4.5, 4.9, 0.9



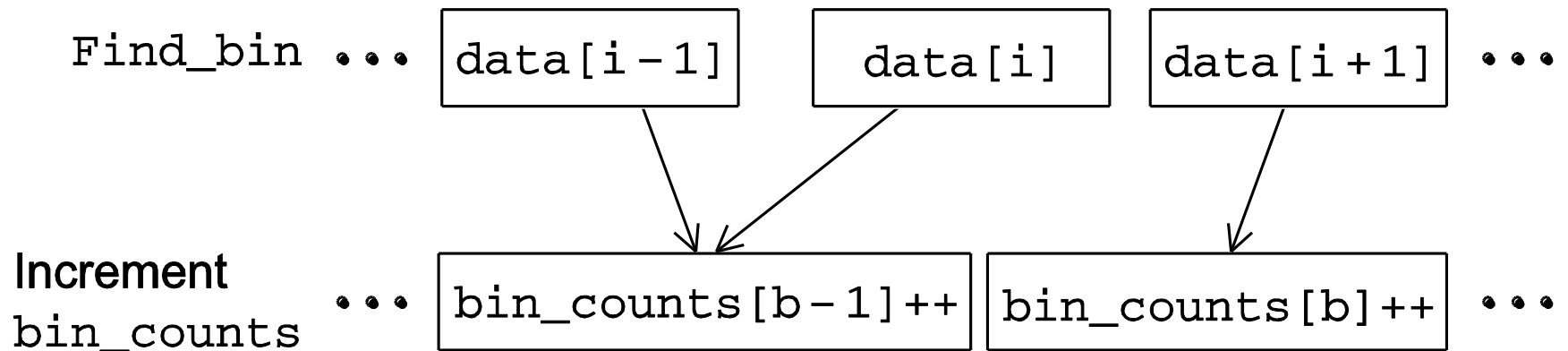
Serial program - input

1. The number of measurements: `data_count`
2. An array of `data_count` floats: `data`
3. The minimum value for the bin containing the smallest values: `min_meas`
4. The maximum value for the bin containing the largest values: `max_meas`
5. The number of bins: `bin_count`

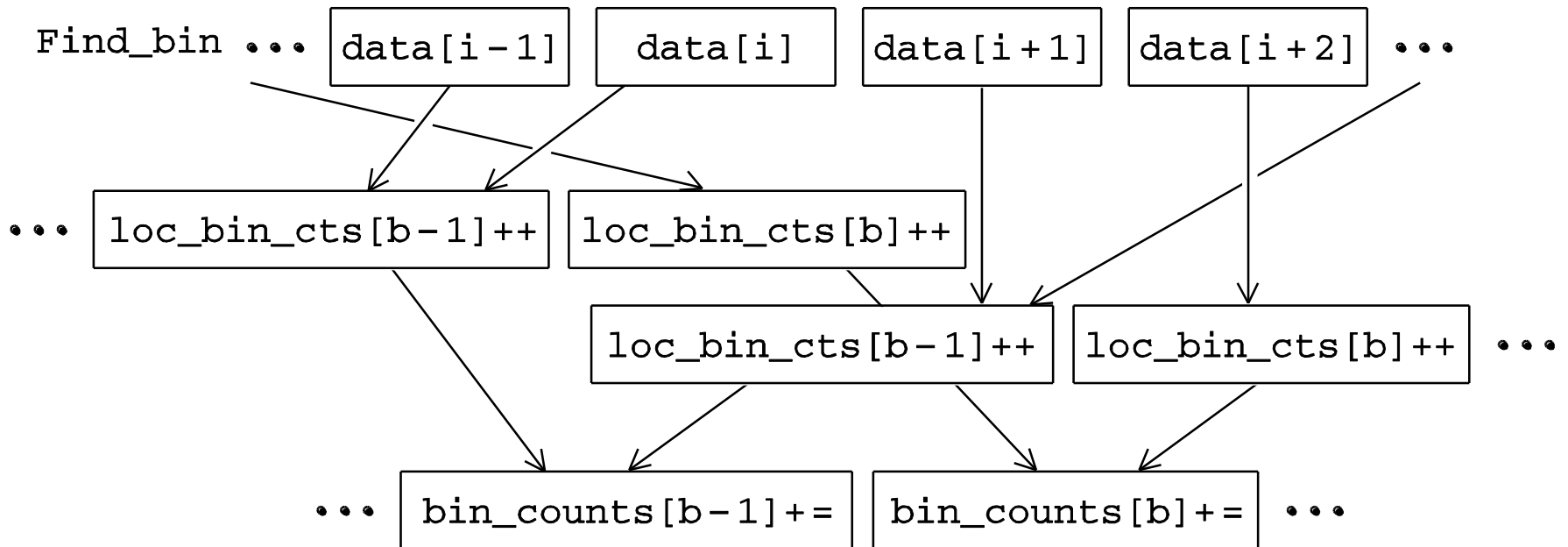
Serial program - output

1. `bin_maxes` : an array of `bin_count` floats
2. `bin_counts` : an array of `bin_count` ints

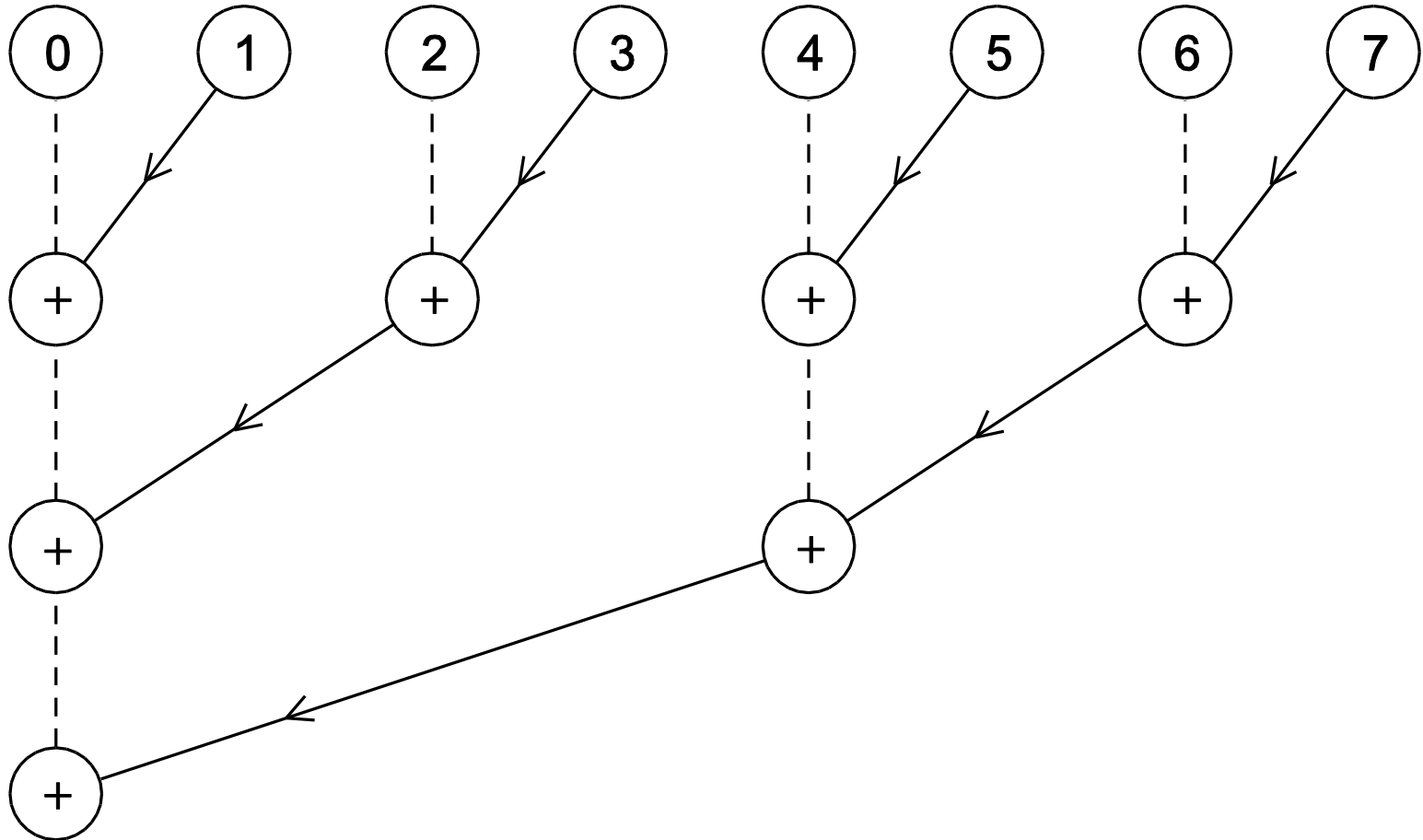
First two stages of Foster's Methodology



Alternative definition of tasks and communication



Adding the local arrays



Concluding Remarks (1)

- Serial systems
 - The standard model of computer hardware has been the von Neumann architecture.
- Parallel hardware
 - Flynn's taxonomy.
- Parallel software
 - We focus on software for homogeneous MIMD systems, consisting of a single program that obtains parallelism by branching.
 - SPMD programs.

Concluding Remarks (2)

- Input and Output
 - We'll write programs in which one process or thread can access stdin, and all processes can access stdout and stderr.
 - However, because of nondeterminism, except for debug output we'll usually have a single process or thread accessing stdout.

Concluding Remarks (3)

- Performance
 - Speedup
 - Efficiency
 - Amdahl's law
 - Scalability
- Parallel Program Design
 - Foster's methodology