



## 4.4 Lecture Summary

## **4 Dataflow Synchronization and Pipelining**

## 4.4 Pipeline Parallelism

**Lecture Summary:** In this lecture, we studied how point-to-point synchronization can be used to build a one-dimensional *pipeline* with p tasks (stages),  $T_0, \ldots, T_{p-1}$ . For example, three important stages in a *medical imaging* pipeline are *denoising*, *registration*, and *segmentation*.

We performed a simplified analysis of the *WORK* and *SPAN* for pipeline parallelism as follows. Let n be the number of input items and p the number of stages in the pipeline,  $WORK = n \times p$  is the total work that must be done for all data items, and CPL = n + p - 1 is the *span* or *critical path length* for the pipeline. Thus, the ideal parallelism is PAR = WORK / CPL = np / (n + p - 1). This formula can be validated by considering a few boundary cases. When p = 1, the ideal parallelism degenerates to PAR = 1, which confirms that the computation is sequential when only one stage is available. Likewise, when n = 1, the ideal parallelism again degenerates to PAR = 1, which confirms that the computation is sequential when only one data item is available. When n is much larger than  $p(n \cdot p)$ , then the ideal parallelism approaches PAR = p in the limit, which is the best possible case.

The synchronization required for pipeline parallelism can be implemented using phasers by allocating an array of phasers, such that phaser  $\mathbf{ph}[\mathbf{i}]$  is "signalled" in iteration i by a call to  $\mathbf{ph}[\mathbf{i}]$ .  $\mathbf{arrive}()$  as follows:

```
1 // Code for pipeline stage i
2 while ( there is an input to be processed ) {
3    // wait for previous stage, if any
4    if (i > 0) ph[i - 1].awaitAdvance();
5
6    process input;
7
8    // signal next stage
9    ph[i].arrive();
10 }
```

## **Optional Reading:**

1. Wikipedia article on <u>Pipeline (computing)</u>.