

# Pipelined Computations

## 1 Functional Decomposition

- car manufacturing with three plants
- speedup for  $n$  inputs in a  $p$ -stage pipeline
- loop unrolling

## 2 Pipeline Implementations

- processors in a ring topology
- pipelined addition
- pipelined addition with MPI

MCS 572 Lecture 25  
Introduction to Supercomputing  
Jan Verschelde, 23 October 2024

# Pipelined Computations

## 1 Functional Decomposition

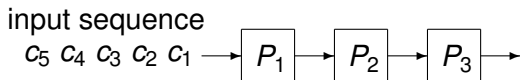
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## 2 Pipeline Implementations

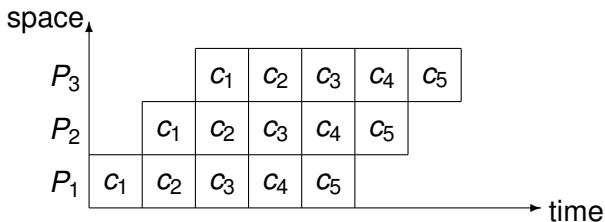
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# car manufacturing

Consider a simplified car manufacturing process in three stages: (1) assemble exterior, (2) fix interior, and (3) paint and finish:



The corresponding *space-time diagram* is below:

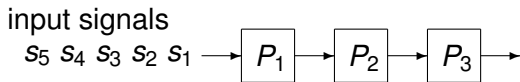


After 3 time units, one car per time unit is completed.

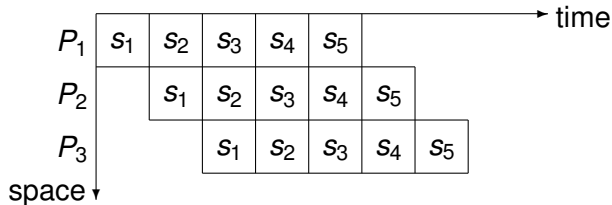
# denoising a signal

Every second we take 256 samples of a signal:

$P_1$ : apply FFT,  $P_2$ : remove low amplitudes, and  $P_3$ : inverse FFT.



An alternative space-time diagram is below:



Observe: the consumption of a signal is sequential.

## $p$ -stage pipelines

A pipeline with  $p$  processors is a  *$p$ -stage pipeline*.

Suppose every process takes one time unit to complete.

How long till a  $p$ -stage pipeline completes  $n$  inputs?

A  $p$ -stage pipeline on  $n$  inputs:

- After  $p$  time units the first input is done.
- Then, for the remaining  $n - 1$  items, the pipeline completes at a rate of one item per time unit.

⇒  $p + n - 1$  time units for the  $p$ -stage pipeline to complete  $n$  inputs.

A time unit is called a *pipeline cycle*.

The time taken by the first  $p - 1$  cycles is the *pipeline latency*.

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# speedup

Consider  $n$  inputs for a  $p$ -stage pipeline:

$$S(p) = \frac{n \times p}{p + n - 1}.$$

For fixed number  $p$  of processors:

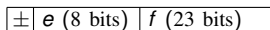
$$\lim_{n \rightarrow \infty} \frac{p \times n}{n + p - 1} = p.$$

Pipelining speeds up multiple sequences of heterogeneous jobs.

- Pipelining is a functional decomposition method to develop parallel programs.
- Recall the classification of Flynn:  
MISD = Multiple Instruction Single Data stream.

# floating-point addition

A floating-point number consists of a sign bit,  
an exponent and a fraction (or mantissa):



Floating-point addition could be done in 6 cycles:

- 1 unpack fractions and exponents
- 2 compare exponents
- 3 align fractions
- 4 add fractions
- 5 normalize result
- 6 pack fraction and exponent of result

Adding two vectors of  $n$  floats with 6-stage pipeline  
takes  $n + 6 - 1$  pipeline cycles, instead of  $6n$  cycles.  
 $\Rightarrow$  Capable of performing one flop per clock cycle.



# Intel Architecture Software Developer's Manual

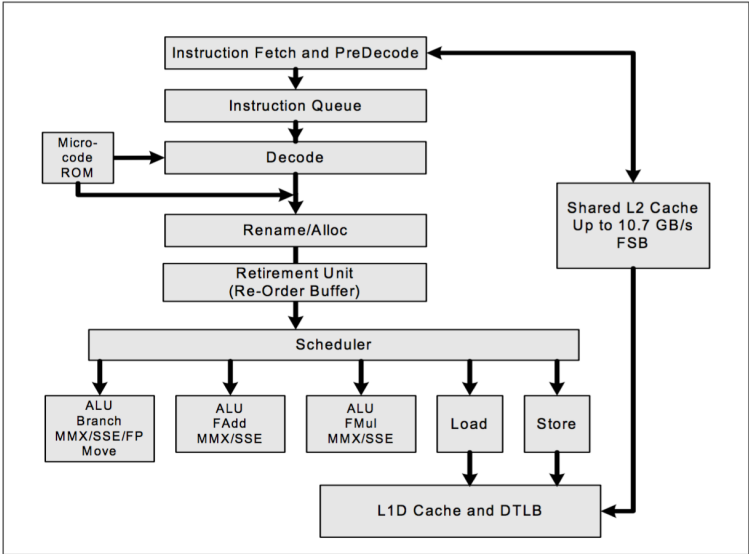


Figure 2-3. The Intel Core Microarchitecture Pipeline Functionality

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# the Leibniz series

The Leibniz series

$$\frac{\pi}{4} = 1 - \frac{1}{3} + \frac{1}{5} - \frac{1}{7} + \frac{1}{9} - \frac{1}{11} + \dots$$

converges very slowly.

This example is based on section 3.2.2 on loop unrolling in *Scientific Programming and Computer Architecture* by Divakar Viswanath, Springer-Verlag, 2017.

The above reference offers a very detailed explanation.

We can already illustrate the main point in Julia.

## a straightforward implementation

The branching in the straightforward code below prevents a pipelined execution of the floating-point operations.

```
function leibniz1(N::Int)
    s = 1.0
    for i=1:N
        if(i%2 == 1)
            s = s - 1.0/(2.0*i + 1.0)
        else
            s = s + 1.0/(2.0*i + 1.0)
        end
    end
    return s
end
```

$$\frac{\pi}{4} = 1 - \frac{1}{3} + \frac{1}{5} - \frac{1}{7} + \frac{1}{9} - \frac{1}{11} + \dots$$

## applying loop unrolling

Summing the even and odd terms separately avoids branching, allows a pipelined executions of the floating-point operations.

```
function leibniz2(N::Int)
    s = 1.0
    for i=2:2:N
        s = s + 1.0/(2.0*i + 1.0)
    end
    for i=1:2:N
        s = s - 1.0/(2.0*i + 1.0)
    end
    return s
end
```

$$1 - \frac{1}{3} + \frac{1}{5} - \frac{1}{7} + \frac{1}{9} - \frac{1}{11} + \dots = 1 + \frac{1}{5} + \frac{1}{9} + \dots - \frac{1}{3} - \frac{1}{7} - \frac{1}{11} - \dots$$

# benchmarking

using BenchmarkTools

```
println(4.0*leibniz1(10^8))  
@btime leibniz1(10^8)
```

```
println(4.0*leibniz2(10^8))  
@btime leibniz2(10^8)
```

with output:

```
3.141592663589326  
 239.600 ms (0 allocations: 0 bytes)  
3.1415926635801443  
 125.266 ms (0 allocations: 0 bytes)
```

Julia 1.8.5 on pascal:

- two 22-core Intel Xeon E5-2699v4 Broadwell at 2.20GHz,
- 256GB of internal memory at 2400MHz.

# Pipelined Computations

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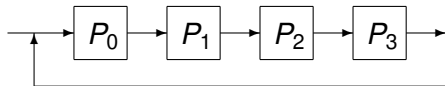
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- processors in a ring topology
- pipelined addition
- pipelined addition with MPI

# processors in a ring topology

A ring topology is a natural way to implement a pipeline.



A manager/worker organization:

- Node 0 receives input and sends to node 1.
- Every node  $i$ , for  $i = 1, 2, \dots, p - 1$ :
  - 1 receives an item from node  $i - 1$ ,
  - 2 performs operations on the item,
  - 3 sends processed item to node  $(i + 1) \bmod p$ .

At the end of one cycle, node 0 has the output.



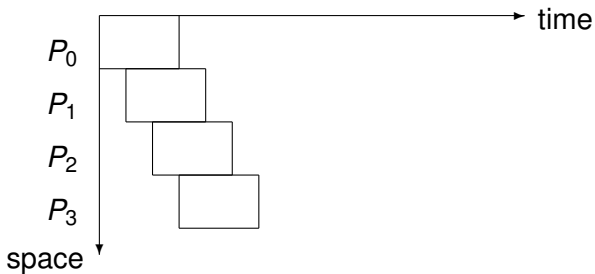
## one pipeline cycle with MPI

```
$ mpirun -np 4 ./pipe_ring
One pipeline cycle for repeated doubling.
Reading a number...
2
Node 0 sends 2 to the pipe...
Processor 1 receives 2 from node 0.
Processor 2 receives 4 from node 1.
Processor 3 receives 8 from node 2.
Node 0 received 16.
$
```

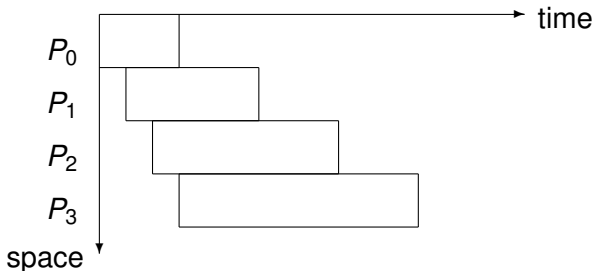
This example is a *type 1 pipeline* efficient only if we have more than one instance to compute.

# space time diagrams for type 2 and type 3 pipelines

Type 2:



Type 3:



## MPI code for the manager

```
void manager ( int p )
/*
 * The manager prompts the user for a number
 * and passes this number to node 1 for doubling.
 * The manager receives from node p-1 the result. */
{
    int n;
    MPI_Status status;

    printf("One pipeline cycle for repeated doubling.\n");
    printf("Reading a number...\n"); scanf("%d",&n);
    printf("Node 0 sends %d to the pipe...\n",n);
    fflush(stdout);
    MPI_Send(&n,1,MPI_INT,1,tag,MPI_COMM_WORLD);
    MPI_Recv(&n,1,MPI_INT,p-1,tag,MPI_COMM_WORLD,&status);
    printf("Node 0 received %d.\n",n);
}
```

## MPI code for the workers

```
void worker ( int p, int i )
/*
 * Worker with identification label i of p
 * receives a number,
 * doubles it and sends it to node i+1 mod p. */
{
    int n;
    MPI_Status status;

    MPI_Recv(&n,1,MPI_INT,i-1,tag,MPI_COMM_WORLD,&status);
    printf("Processor %d receives %d from node %d.\n",
           i,n,i-1);
    fflush(stdout);
    n *= 2;                               /* double the number */
    if(i < p-1)
        MPI_Send(&n,1,MPI_INT,i+1,tag,MPI_COMM_WORLD);
    else
        MPI_Send(&n,1,MPI_INT,0,tag,MPI_COMM_WORLD);
}
```

# Pipelined Computations

## 1 Functional Decomposition

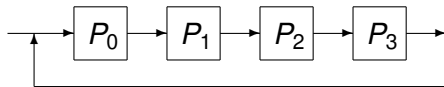
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## 2 Pipeline Implementations

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- **pipelined addition**
- pipelined addition with MPI

# pipelined addition

Consider 4 processors in a ring topology:



To add a sequence of 32 numbers, with data partitioning:

$$\underbrace{a_0, a_1, \dots, a_7}, \underbrace{b_0, b_1, \dots, b_7}, \underbrace{c_0, c_1, \dots, c_7}, \underbrace{d_0, d_1, \dots, d_7}.$$
$$A_k = \sum_{j=0}^k a_j \quad B_k = \sum_{j=0}^k b_j \quad C_k = \sum_{j=0}^k c_j \quad D_k = \sum_{j=0}^k d_j$$

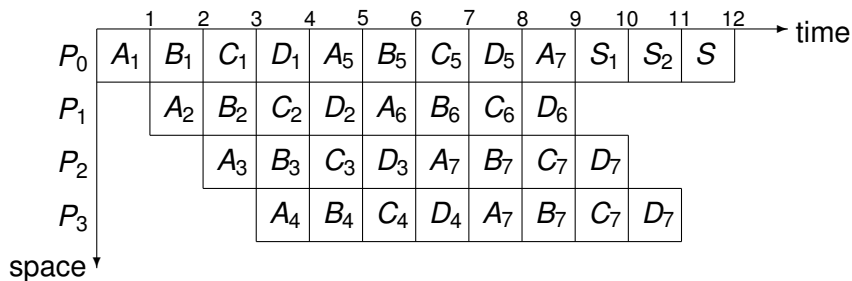
The final sum is  $S = A_7 + B_7 + C_7 + D_7$ .

# space-time diagram for pipelined addition

$$\underbrace{a_0, a_1, \dots, a_7}, \underbrace{b_0, b_1, \dots, b_7}, \underbrace{c_0, c_1, \dots, c_7}, \underbrace{d_0, d_1, \dots, d_7}.$$

$$A_k = \sum_{j=0}^k a_j \quad B_k = \sum_{j=0}^k b_j \quad C_k = \sum_{j=0}^k c_j \quad D_k = \sum_{j=0}^k d_j$$

Denote  $S_1 = A_7 + B_7$ ,  $S_2 = S_1 + C_7$ ,  $S = S_2 + D_7$ .



## speedup for pipelined addition

We finished addition of 32 numbers in 12 cycles:  $12 = 32/4 + 4$ .

In general, with  $p$ -stage pipeline to add  $n$  numbers:

$$S(p) = \frac{n-1}{\frac{n}{p} + p}$$

For fixed  $p$ :  $\lim_{n \rightarrow \infty} S(p) = p$ .



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## using 5-stage pipeline

```
$ mpirun -np 5 ./pipe_sum
```

```
The data to sum : 1 2 3 4 5 6 7 8 9 10 11 12 13 14 \
15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30
```

```
Manager starts pipeline for sequence 0...
```

```
Processor 1 receives sequence 0 : 3 3 4 5 6
```

```
Processor 2 receives sequence 0 : 6 4 5 6
```

```
Processor 3 receives sequence 0 : 10 5 6
```

```
Processor 4 receives sequence 0 : 15 6
```

```
Manager received sum 21.
```

```
Manager starts pipeline for sequence 1...
```

```
Processor 1 receives sequence 1 : 15 9 10 11 12
```

```
Processor 2 receives sequence 1 : 24 10 11 12
```

```
Processor 3 receives sequence 1 : 34 11 12
```

```
Processor 4 receives sequence 1 : 45 12
```

```
Manager received sum 57.
```

## session continued

Manager starts pipeline for sequence 2...

Processor 1 receives sequence 2 : 27 15 16 17 18

Processor 2 receives sequence 2 : 42 16 17 18

Processor 3 receives sequence 2 : 58 17 18

Processor 4 receives sequence 2 : 75 18

Manager received sum 93.

Manager starts pipeline for sequence 3...

Processor 1 receives sequence 3 : 39 21 22 23 24

Processor 2 receives sequence 3 : 60 22 23 24

Processor 3 receives sequence 3 : 82 23 24

Processor 4 receives sequence 3 : 105 24

Manager received sum 129.

## end of the session

```
Manager starts pipeline for sequence 4...
Processor 1 receives sequence 4 : 51 27 28 29 30
Processor 2 receives sequence 4 : 78 28 29 30
Processor 3 receives sequence 4 : 106 29 30
Processor 4 receives sequence 4 : 135 30
Manager received sum 165.
The total sum : 465
$
```

## MPI code

```
void pipeline_sum ( int i, int p )
/* performs a pipeline sum of p*(p+1) numbers */
{
    int n[p][p-i+1];
    int j,k;
    MPI_Status status;

    if(i==0) /* manager generates numbers */
    {
        for(j=0; j<p; j++)
            for(k=0; k<p+1; k++) n[j][k] = (p+1)*j+k+1;
        if(v>0)
        {
            printf("The data to sum : ");
            for(j=0; j<p; j++)
                for(k=0; k<p+1; k++) printf(" %d",n[j][k]);
            printf("\n");
        }
    }
}
```

# loop for manager

```
for(j=0; j<p; j++)
  if(i==0) /* manager starts pipeline of j-th sequence */
  {
    n[j][1] += n[j][0];
    printf("Manager starts pipeline for sequence %d...\n",
           j);
    MPI_Send(&n[j][1], p, MPI_INT, 1, tag, MPI_COMM_WORLD);
    MPI_Recv(&n[j][0], 1, MPI_INT, p-1, tag, MPI_COMM_WORLD,
             &status);
    printf("Manager received sum %d.\n", n[j][0]);
  }
else /* worker i receives p-i+1 numbers */
```

## loop for workers

```
else      /* worker i receives p-i+1 numbers */
{
    MPI_Recv(&n[j][0],p-i+1,MPI_INT,i-1,tag,
             MPI_COMM_WORLD,&status);
    printf("Processor %d receives sequence %d : ",i,j);
    for(k=0; k<p-i+1; k++) printf(" %d", n[j][k]);
    printf("\n");
    n[j][1] += n[j][0];
    if(i < p-1)
        MPI_Send(&n[j][1],p-i,MPI_INT,i+1,tag,
                 MPI_COMM_WORLD);
    else
        MPI_Send(&n[j][1],1,MPI_INT,0,tag,MPI_COMM_WORLD);
}
if(i==0) /* manager computes the total sum */
{
    for(j=1; j<p; j++) n[0][0] += n[j][0];
    printf("The total sum : %d\n",n[0][0]);
}
```

# Summary + Exercises

We started chapter 5 in the book of Wilkinson and Allen.

## Exercises:

- 1 Describe the application of pipelining technique for grading  $n$  copies of an exam that has  $p$  questions. Explain the stages and make a space-time diagram.
- 2 Write code to use the 4-stage pipeline to double numbers for a sequence of 10 consecutive numbers starting at 2.
- 3 Consider the evaluation of a polynomial  $f(x)$  of degree  $d$  given by its coefficient vector  $(a_0, a_1, a_2, \dots, a_d)$ , using Horner's method, e.g., for  $d = 4$ :  $f(x) = (((a_4x + a_3)x + a_2)x + a_1)x + a_0$ .  
Give MPI code of this algorithm to evaluate  $f$  at a sequence of  $n$  values for  $x$  by a  $p$ -stage pipeline.