#### **EH**zürich

TIMO SCHNEIDER (SUBST. TORSTEN HOEFLER)

# Parallel Programming Finish STM & Distributed Memory Programming: Actors, CSP, and MPI

#### Microsoft, Google: We've found a fourth data-leaking Meltdown-Spectre CPU hole

Design blunder exists in Intel, AMD, Arm, Power processors

By Chris Williams, Editor in Chie 21 May 2018 at 21:00 74 📮 SHARE 🔻



A fourth variant of the data-leaking Meltdown-Spectre security flaws in modern processors has been found by Microsoft and Google researchers.



How the fourth variant works

Variant 4 is referred to as a speculative store bypass. It is yet another "wait, why didn't I think of that?" design oversight in modern out-of-orderexecution engineering. And it was found by Google Project Zero's Jann Horn, who helped uncover the earlier Spectre and Meltdown bugs, and Ken Johnson of Microsoft.

It hinges on the fact that when faced with a bunch of software instructions that store data to memory, the CPU will look far ahead to see if it can execute any other instructions out of order while the stores complete. Writing to memory is generally slow compared to other instructions. A modern fast CPU won't want to be held up by store operations, so it looks ahead to find other things to do in the meantime.

If the processor core, while looking ahead in a program, finds an instruction that loads data from memory, it will predict whether or not this load operation is affected by any of the preceding stores. For example, if a store is writing to memory that a later load fetches back from memory, you'll want the store to complete first. If a load is predicted to be safe to run ahead of the pending stores, the processor executes it speculatively while other parts of the chip are busy with other code.



#### Bank account (ScalaSTM)

```
class AccountSTM {
    private final Integer id; // account id
    private final Ref.View<Integer> balance;
```

```
AccountSTM(int id, int balance) {
   this.id = new Integer(id);
   this.balance = STM.newRef(balance);
```



#### Ideal world: bank account using atomic keyword

```
void withdraw(final int amount) {
     // assume that there are always sufficient funds...
     atomic {
          int old_val = balance.get();
          balance.set(old val - amount);
      }
void deposit(final int amount) {
     atomic {
          int old val = balance.get();
          balance.set(old val + amount);
      }
```



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#### Real world: bank account in ScalaSTM

```
void withdraw(final int amount) {
     // assume that there are always sufficient funds...
     STM.atomic(new Runnable() { public void run() {
         int old val = balance.get();
         balance.set(old val - amount);
     }});
void deposit(final int amount) {
     STM.atomic(new Runnable() { public void run() {
         int old val = balance.get();
         balance.set(old val + amount);
     }});
```



#### **GetBalance (return a value)**

```
public int getBalance() {
  int result = STM.atomic(
    new Callable<Integer>() {
    public Integer call() {
"atomic"
      int result = balance.get();
      return result;
  });
  return result;
}
```



#### **Bank account transfer**

What if account a does not have enough funds?

How can we wait until it does in order to retry the transfer?

#### locks $\rightarrow$ conditional variables

 $TM \rightarrow retry$ 

#### Bank account transfer with retry

```
static void transfer_retry(final AccountSTM a,
                            final AccountSTM b,
                            final int amount) {
     atomic {
           if (a.balance.get() < amount)</pre>
                STM.retry();
           a.withdraw(amount);
           b.deposit(amount);
```

retry: abort the transaction and retry when conditions change



### How does retry work?

Implementations need to track what reads/writes a transaction performed to detect conflicts

- Typically called read-/write-set of a transaction
- When retry is called, transaction aborts and will be retried when any of the variables that were read, change
- In our example, when a.balance is updated, the transaction will be retried



#### **Simplest STM Implementation**

# Ingredients

Threads that run transactions with thread states

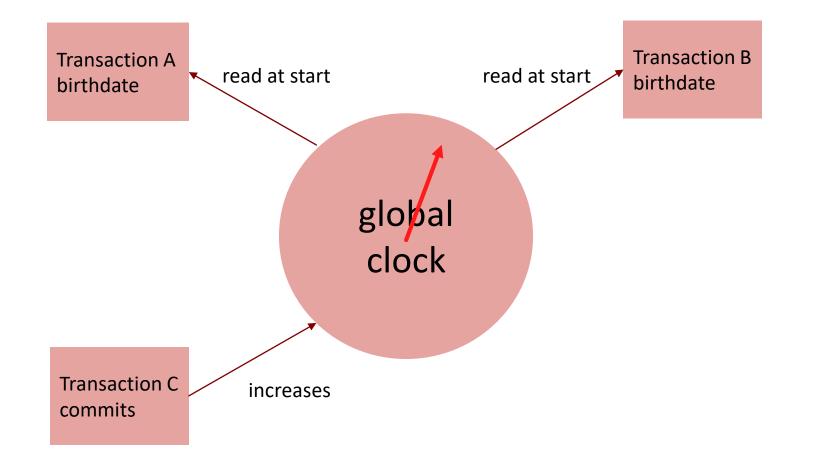
- active
- aborted
- committed

Objects representing state stored in memory (the variables affected by a transaction)

- offering methods like a constructor, read (get), write (set)
- and copy!



#### **Clock-based STM System**



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# **Atomic Objects**

Each transaction uses a local **read-set** and a local **write-set** holding all locally read and written objects.

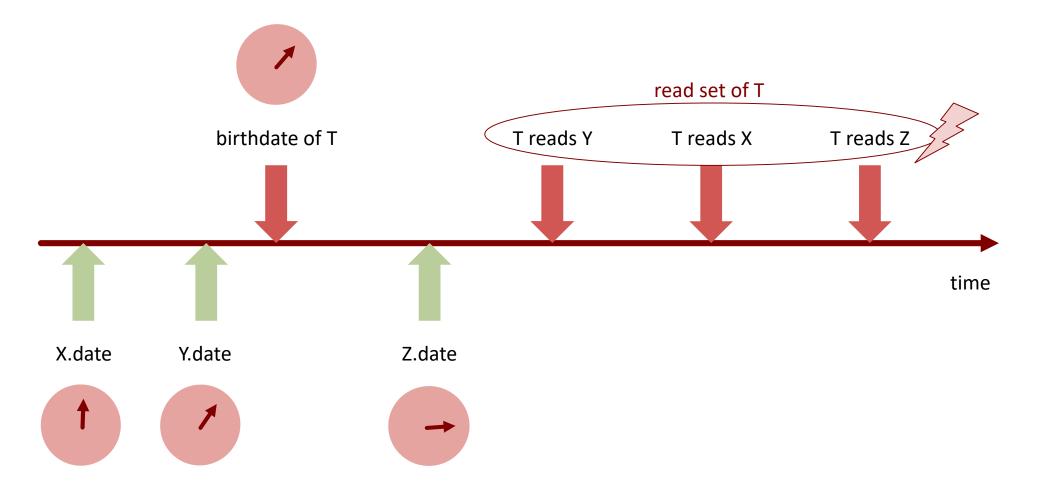
Transaction calls read

atomic memory object version time reference stamp

- check if the object is in the write set ightarrow return this (new) version
- otherwise check if object's time stamp ≤ transaction's birthdate, if not throw aborted exception, otherwise add new copy of the object to the read set
- Transaction calls write
- if object is not in write set, create a copy of it in the write set



#### **Transaction life time**



Martin Come

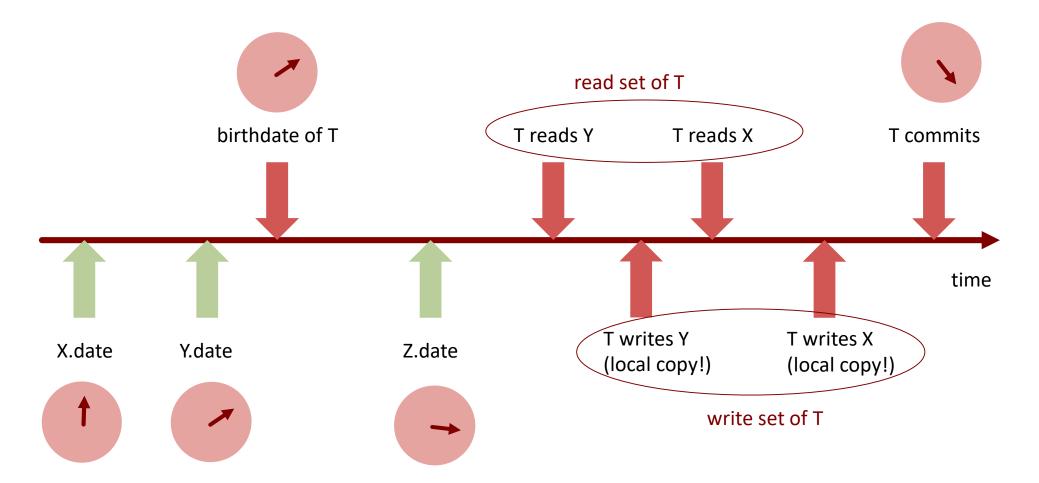


# Commit

- Lock all objects of read- and write-set (in some defined order to avoid deadlocks)
- Check that all objects in the read set provide a time stamp ≤ birthdate of the transaction, otherwise return "abort"
- Increment and get the value T of current global clock
- Copy each element of the write set back to global memory with timestamp T
- Release all locks and return "commit"



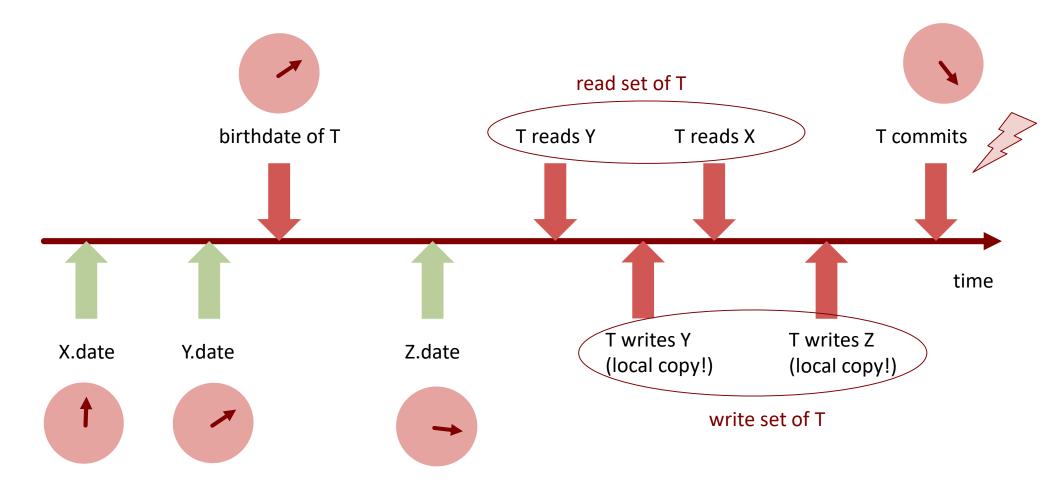
#### Successful commit



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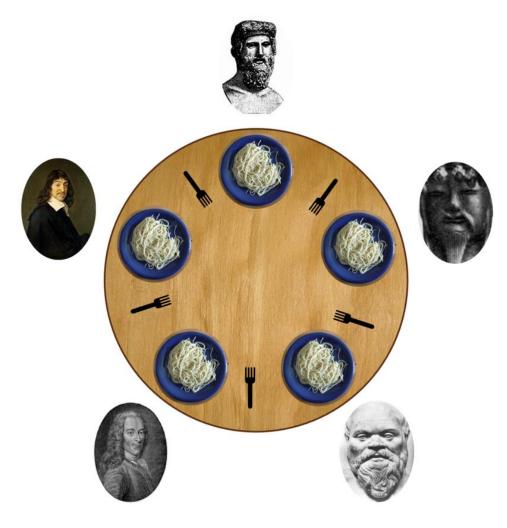
#### **Aborted commit**



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# **Dining philosophers**



- 5 philosophers
- 5 forks
- each philosopher requires 2 forks to eat
- forks cannot be shared

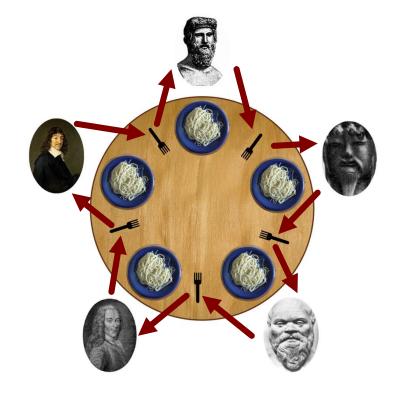
image source: Wikipedia



### Solution that can lead to deadlock

Philosopher:

- think
- lock left
- lock right
- eat
- unlock right
- unlock left



 $P_1$  takes  $F_1$ ,  $P_2$  takes  $F_2$ ,  $P_3$  takes  $F_3$ ,  $P_4$  takes  $F_4$ ,  $P_5$  takes  $F_5$  $\rightarrow$  Deadlock



#### **Dining Philosophers Using TM**

```
private static class Fork {
     public final Ref.View<Boolean> inUse = STM.newRef(false);
}
class PhilosopherThread extends Thread {
      private final int meals;
      private final Fork left;
     private final Fork right;
      public PhilosopherThread(Fork left, Fork right) {
           this.left = left;
           this.right = right;
      public void run() { ... }
```

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#### **Dining Philosophers Using TM**

```
Fork[] forks = new Fork[tableSize];
```

```
for (int i = 0; i < tableSize; i++)
    forks[i] = new Fork();</pre>
```

PhilosopherThread[] threads = new PhilosopherThread[tableSize];

A State of the owned

```
for (int i = 0; i < tableSize; i++)
    threads[i] = new PhilosopherThread(forks[i],
        forks[(i + 1) % tableSize]);</pre>
```



Contra anna anna anna

### **Dining Philosophers Using TM**

```
class PhilosopherThread extends Thread {
   ...
      public void run() {
            for (int m = 0; m < meals; m++) {</pre>
                  // THINK
                  pickUpBothForks();
                  // EAT
                  putDownForks();
            }
   ...
```



...

...

```
Dining Philosophers Using TM
```

class PhilosopherThread extends Thread {

```
private void pickUpBothForks() {
   STM.atomic(new Runnable() { public void run() {
```

```
if (left.inUse.get() || right.inUse.get())
    STM.retry();
```

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```
left.inUse.set(true);
right.inUse.set(true);
```

```
}});
```



...

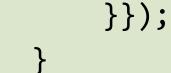
#### **Dining Philosophers Using TM**

class PhilosopherThread extends Thread {

```
private void putDownForks() {
   STM.atomic(new Runnable() { public void run() {
```

```
left.inUse.set(false);
right.inUse.set(false);
```

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#### **Issues with transactions**

- It is not clear what are the best semantics for transactions
- Getting good performance can be challenging
- I/O operations (e.g., print to screen)
   Can we perform I/O operations in a transaction?



#### Summary

- Locks are too hard!
- Transactional Memory tries to remove the burden from the programmer
  STM / HTM
- Remains to be seen whether it will be widely adopted in the future



# **Additional Reading**

Simon Peyton Jones, Beautiful concurrency http://research.microsoft.com/pubs/74063/beautiful.pdf

Dan Grossman,

The Transactional Memory / Garbage Collection Analogy https://homes.cs.washington.edu/~djg/papers/analogy\_oopsla07.pdf



# Distributed Memory & Message Passing

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# So far

### Considered

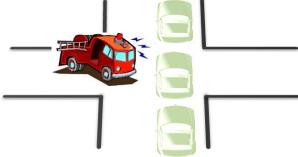
- Parallel / Concurrent
- Fork-Join / Threads
- OOP on Shared Memory
- Locking / Lock Free / Transactional
- Semaphores / Monitors



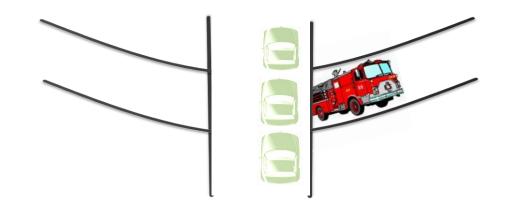
# **Sharing State**

Many of the problems of parallel/concurrent programming come from sharing state

Complexity of locks, race conditions, ....



What if we avoid sharing state?





### Alternatives

# **Functional Programming**

• Immutable state  $\rightarrow$  no synchronization required

## Message Passing: Isolated mutable state

- State is mutable, but not shared: Each thread/task has its private state
- Tasks cooperate via message passing



#### **Concurrent Message Passing**

**Programming Models** 

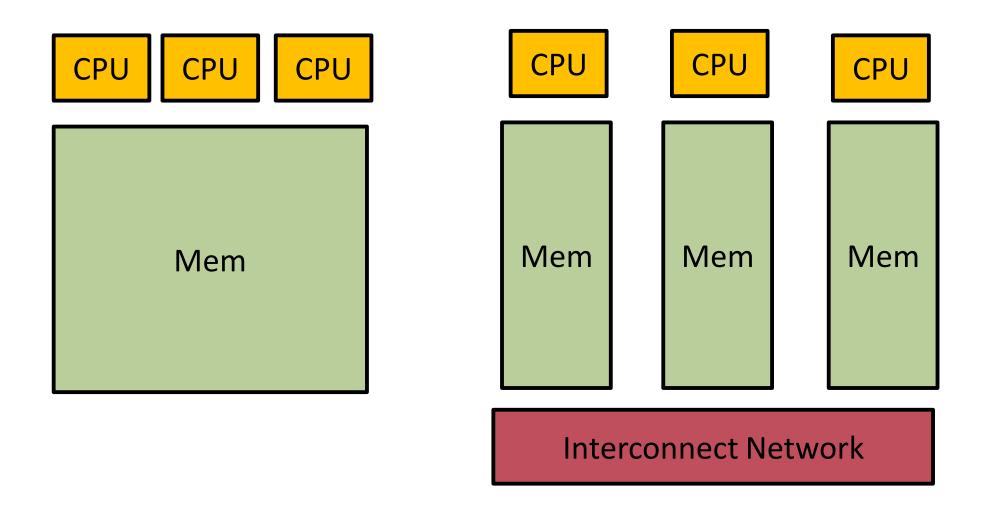
- CSP: Communicating Sequential Processes
- Actor programming model

Framework/library

MPI (Message Passing Interface)



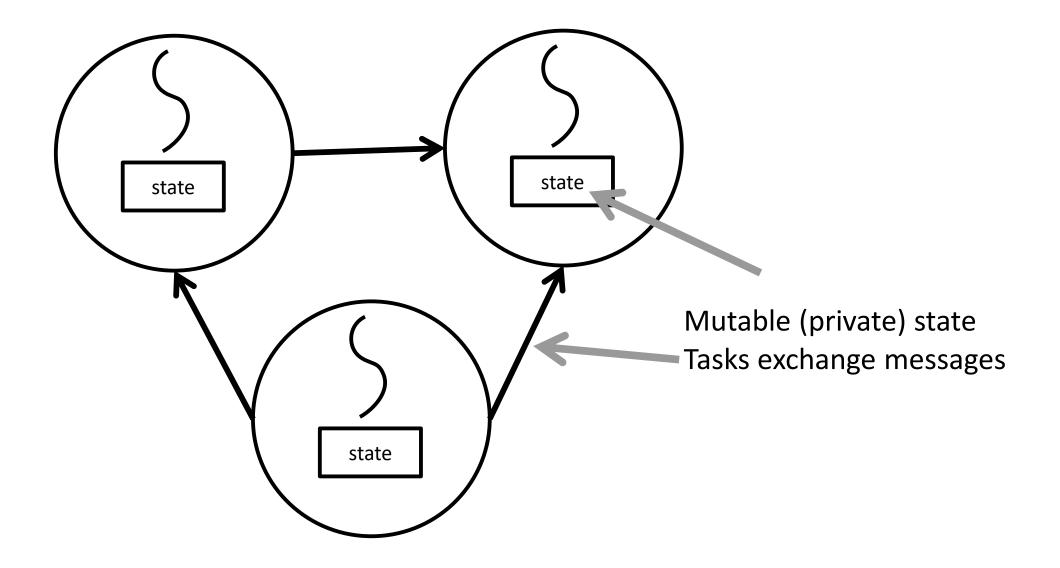
## **Shared vs Distributed memory**



A Real Manager



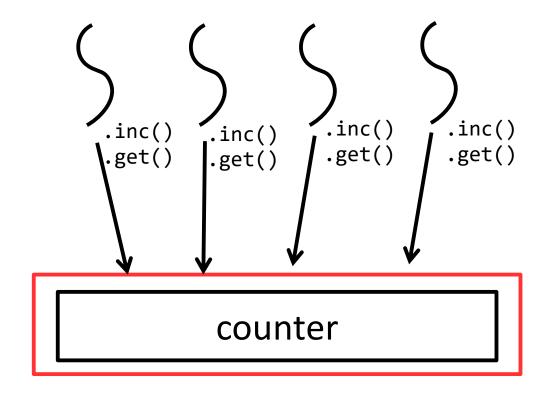
#### **Isolated mutable state**



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#### **Example: Shared state counting**

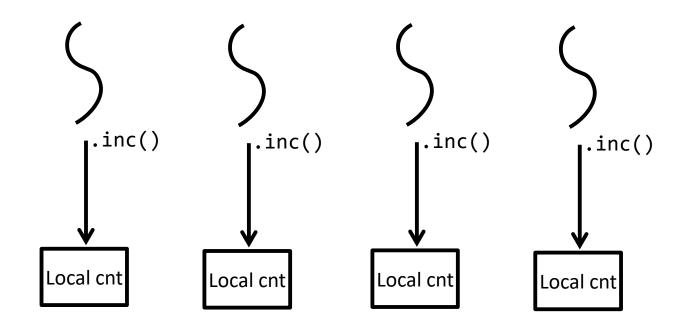


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 $\rightarrow$  shared state must be protected (lock/atomic counter)



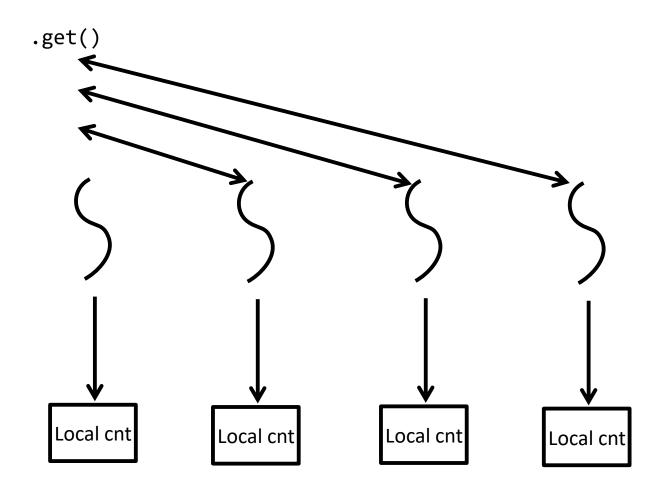
#### **Isolated mutability: counting**



Mal Chandres - 19



**Isolated mutability: accessing count** 



Static Contraction



# **Rethinking managing state**

Bank account

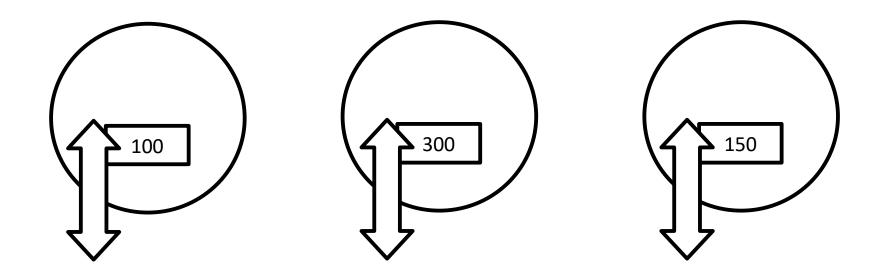
- Sequential programming
  - Single balance
- Parallel programming: sharing state
  - Single balance + protection
- Parallel programming: distributing state
  - Each thread has a local balance (a budget)
  - Threads exchange amounts at coarse granularity



### **Distributed Bank account**

Total balance: 100 + 300 + 150 = 550

- Each task can operate independently
- And communicate with other tasks only when needed
  - This lecture: via messaging





#### Synchronous vs Asynchronous messages

Synchronous:

- sender blocks until message is received



- sender does not block (fire-and-forget)
- placed into a buffer for receiver to get







# The Actor Model\*

Actor = Computational agent that maps

communication to

- a finite set of communications sent to other actors (messages)
- a new behavior (state)
- a finite set of new actors created (dynamic reconfigurability)
- Undefined global ordering
- Asynchronous Message Passing
- Invented by Carl Hewitt 1973\*\*

Actor			
	Thread	Maill	хос
State			



#### **The Actor Model**

Actor model provides a dynamic interconnection topology

- dynamically configure the graph during runtime (add channels)
- dynamically allocate resources

An actor sends messages to other actors using "direct naming", without indirection via port / channel / queue / socket (etc.)

Implemented in various languages such as Erlang, Scala, Ruby and in frameworks such as Akka (for Scala and Java)



#### **Event-driven programming model**

Typically actors react to messages

Event-driven model

A program is written as a set of event handlers for events (events can be seen as received messages)

Example: Graphical User Interface

- user presses OK button  $\rightarrow \dots$
- user presses Cancel button  $\rightarrow \dots$

# **Example: Erlang**

#### Functional Programming Language

- code might look unconventional at first
   Developed by Ericsson for distributed faulttolerant applications
- if no state is shared, recovering from errors becomes much easier

Open source

Concurrent, follows the actor model

```
-module(pingpong).
-export([start/1, ping/2, pong/0]).
ping(0, Pong Node) ->
                                                ERLANG
    {pong, Pong Node} ! finished,
    io:format("ping finished~n", []);
ping(N, Pong Node) ->
   {pong, Pong Node} ! {ping, self()},
    receive
        pong ->
            io:format("Ping received pong~n", [])
    end,
    ping(N - 1, Pong Node).
pong() ->
    receive
        finished \rightarrow
            io:format("Pong finished~n", []);
        {ping, Ping PID} ->
            io:format("Pong received ping~n", []),
            Ping PID ! pong,
            pong()
    end.
start(Ping Node) ->
    register(pong, spawn(pingpong, pong, [])),
    spawn(Ping Node, pingpong, ping, [3, node()]).
```





# **Erlang example**

```
start() ->
      Pid = spawn(fun() -> hello() end),
      Pid ! hello,
      Pid ! bye.
hello() ->
      receive
          hello ->
             io:fwrite("Hello world\n"),
                 hello();
          bye ->
             io:fwrite("Bye cruel world\n"),
                    ok
      end.
```

new task (actor) that will execute the hello function spawn returns address (Pid) of new task

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Address (Pid) can be used to send messages to task

Messages sent to a task are put in a mailbox

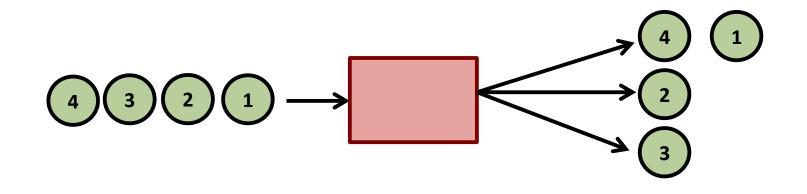
**Receive** reads the first message in the mailbox, which is matched against patterns (similar to a switch statement)

Event-driven programming: code is structured as reactions to events



#### Actor example: distributor

• Forward received messages to a set of nodes in a round-robin fashion





#### Actor example: distributor

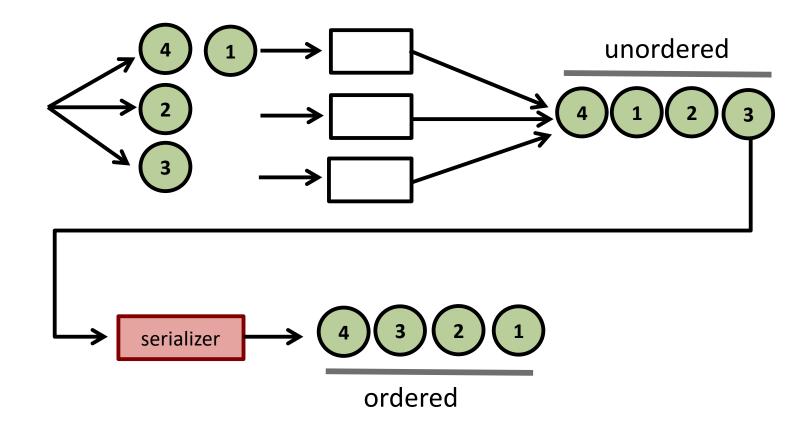
State:

- an array of actors
- the array index of the next actor to forward a message

Receive:

- messages  $\rightarrow$  forward message and increase index (mod)
- control commands (e.g., add/remove actors)



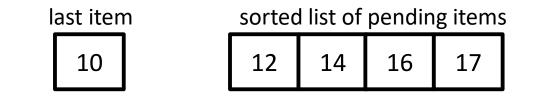


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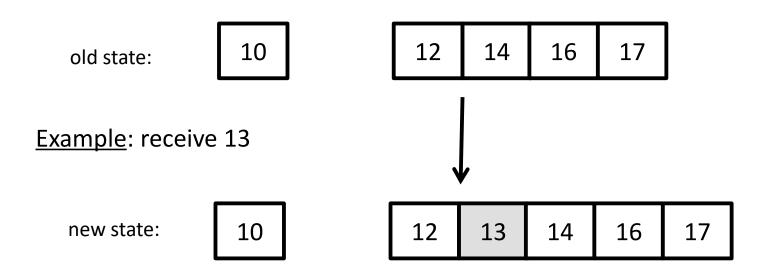
State:

- a sorted list of items we have received
- the last item we forwarded



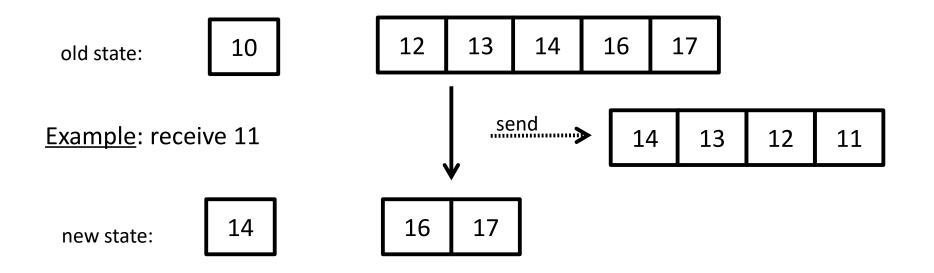


Receive: If we receive an item that is <u>larger</u> than the last item plus one: – add it to the sorted list





- Receive: If we receive an item that is <u>equal to</u> the last item plus one:
  - send the received item plus all consecutive items from the list
  - reset the last item





# Communicating Sequential Processes (1978, 1985)

# Sir Charles Antony Richard Hoare (aka C.A.R. / Tony Hoare)

Formal language defining a process algebra for concurrent systems.



Operators seq (sequential) and par (parallel) for the hierarchical composition of processes.

Synchronisation and Communication between parallel processes with Message Passing.

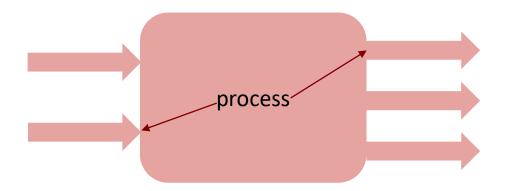
- Symbolic channels between sender and receiver
- Read and write requires a rendezvouz (synchronous!)

CSP was first implemented in Occam.



### **CSP: Indirect Naming**

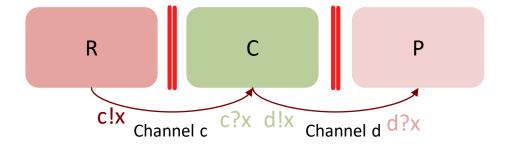
- Many message passing architectures (such as CSP) include an intermediary entity (*port / channel*) to address send destination
- Process issuing send() specifies the port to which the message is sent
- Process issuing receive() specifies a port number and waits for the first message that arrives at the port



# CSP Example (from Hoare's seminal Paper)

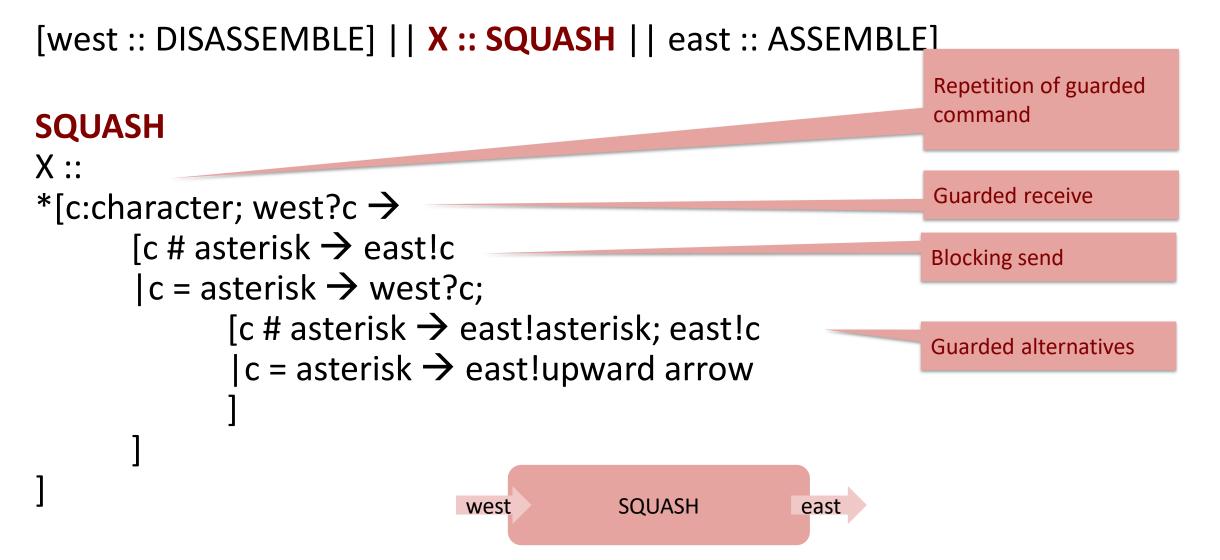
Conway's Problem

- Write a program that transforms a series of cards with 80-character columns in a series of printing lines with 125 characters each. Replace each "\*\*" by "^"
- Separation into processes (Threads)
   R par C par P
  - R: Reading process reading 80-character records
  - C: Converting process converting "\*\*" into "^"
  - P: Printing process: write records with 125 characters





# **CSP Example (from Hoare's seminal Paper)**





#### OCCAM

First programming language to implement CSP (1983)

2 Carlor Maria

```
ALT
 count1 < 100 & c1 ? data
  SEQ
   count1 := count1 + 1
   merged ! data
 count2 < 100 & c2 ? data
  SEQ
   count2 := count2 + 1
   merged ! data
 status ? request
  SEQ
   out ! count1
   out ! count2
```



# Go programming language

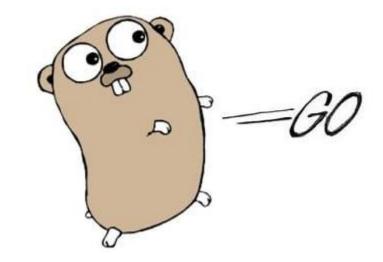
Concurrent programming language from Google

Language support for:

- Lightweight tasks (called goroutines)
- Typed channels for task communications
  - channels are synchronous (or unbuffered) by default
  - support for asynchronous (buffered) channels

Inspired by CSP

Language roots in Algol Family: Pascal, Modula, Oberon [Prof. Niklaus Wirth, ETH] [One of the inventors, Robert Griesemer: PhD from ETH]





}

#### Go example

```
func main() {
       msgs := make(chan string)
       done := make(chan bool)
       go hello(msgs, done);
       msgs <- "Hello"</pre>
       msgs <- "bye"</pre>
       ok := <-done
       fmt.Println("Done:", ok);
```

```
func hello(msgs chan string,
           done chan bool) {
       for {
               msg := <-msgs
               fmt.Println("Got:", msg)
               if msg == "bye" {
                       break
               }
        }
       done <- true;</pre>
```

The second second

}



```
func hello(msgs chan string,
func main() {
                                                               done chan bool) {
        msgs := make(chan string)
        done := make(chan bool)
                                                           for {
                                                                   msg := <-msgs
        go hello(msgs, done);
                                                                   fmt.Println("Got:", msg)
                                   Create two channels:
        msgs <- "Hello"</pre>
                                                                   if msg == "bye" {
                                   •msgs: for strings
                                                                           break
        msgs <- "bye"

    done: for boolean values

                                                                   }
        ok := <-done
        fmt.Println("Done:", ok);
                                                           done <- true;</pre>
                                                   }
```

Manual and the



```
func hello(msgs chan string,
func main() {
                                                               done chan bool) {
       msgs := make(chan string)
       done := make(chan bool)
                                                          for {
                                                                  msg := <-msgs
       go hello(msgs, done);
                                                                  fmt.Println("Got:", msg)
                                 Create a new task (goroutine),
                                                                  if msg == "bye" {
       msgs <- "Hello"
                                 that will execute function
                                                                          break
       msgs <- "bye"</pre>
                                 hello with the given
                                                                  }
                                 arguments
       ok := <-done
       fmt.Println("Done:", ok);
                                                          done <- true;</pre>
                                                  }
```

Contraction of the





```
func main() {
```

```
msgs := make(chan string)
done := make(chan bool)
go hello(msgs, done);
msgs <- "Hello"
msgs <- "bye"</pre>
```

```
ok := <-done
```

```
fmt.Println("Done:", ok);
```

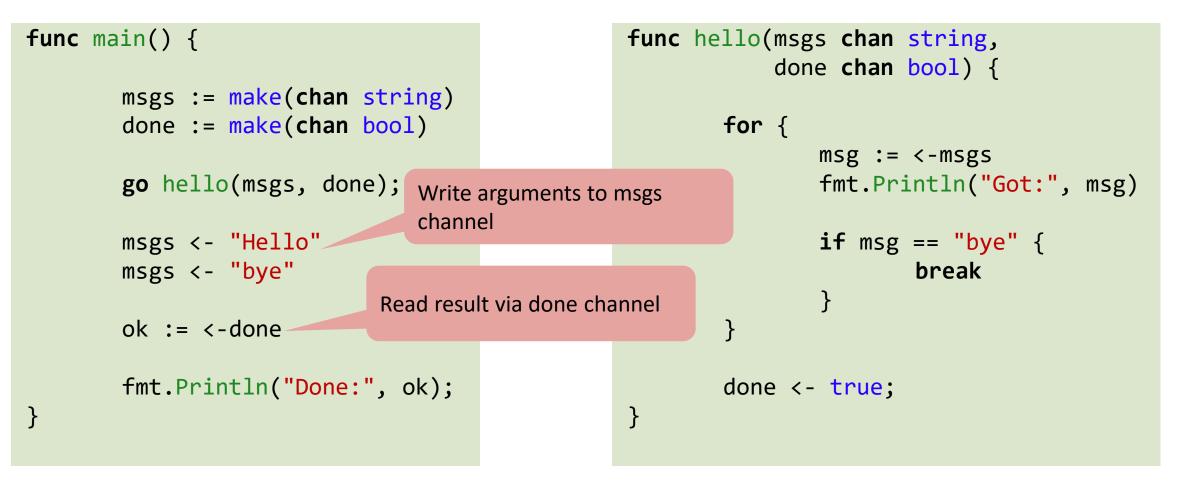
```
Hello takes two channels as
arguments for communication
```

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}

```
func hello(msgs chan string,
           done chan bool) {
       for {
               msg := <-msgs
               fmt.Println("Got:", msg)
               if msg == "bye" {
                       break
               }
        }
       done <- true;</pre>
```





The second state

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### Q: what will happen in this program?

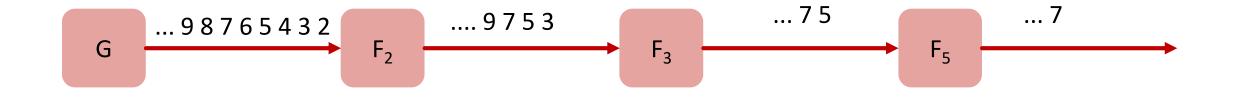
```
func t(in chan string, done chan bool) {
      m := <-in
                        // receive from in channel
      fmt.Println("Got message:", m); // print received message
                     // send true to done channel
      done <- true
}
func main() {
      c := make(chan string) // create a string channel
      done := make(chan bool) // create a boolean channel
      go t(c,done) // spawn goroutine
                 // receive from done channel
      ok := <-done
      fmt.Println("Got ok:", ok); // print ok
      c <- "Hello"
                 // send hello to channel c
```

A: fatal error: all goroutines are asleep - deadlock!



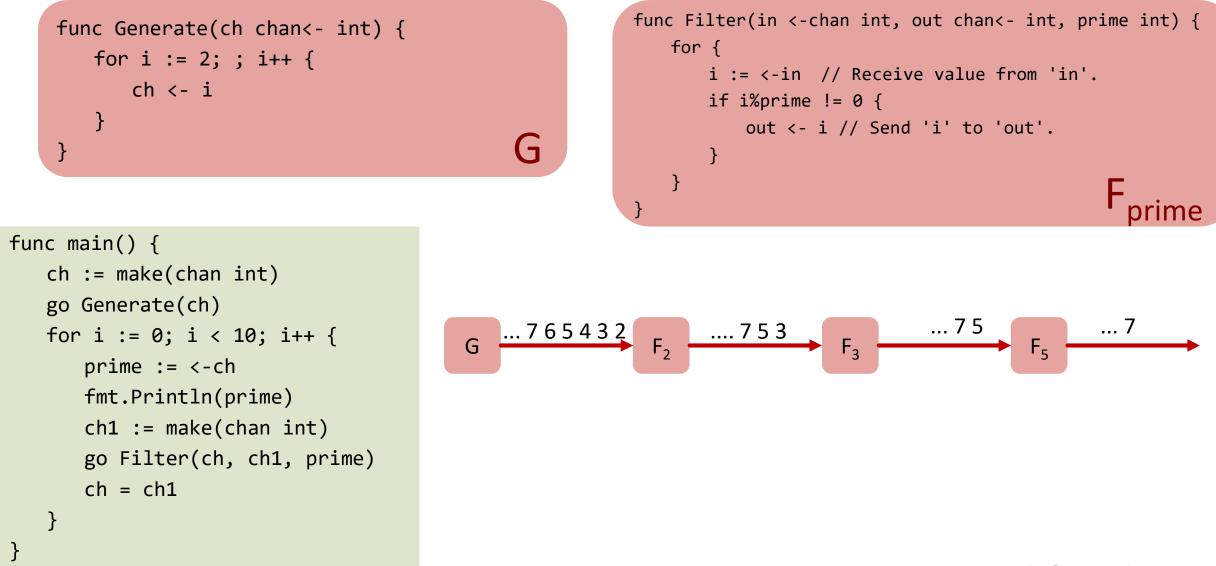
#### **Example: Concurrent prime sieve**

Each station removes multiples of the first element received and passes on the remaining elements to the next station





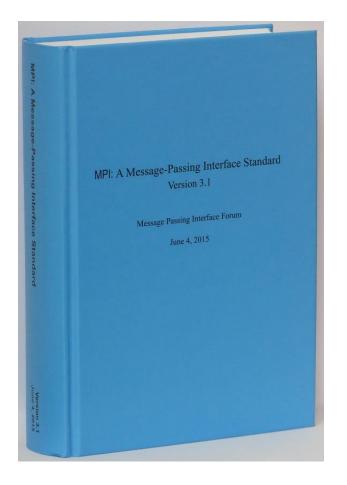
#### **Concurrent prime sieve**



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# Message Passing Interface (MPI)



#### Parallel Processing Letters

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#### MPI ON MILLIONS OF CORES\*

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#### ABSTRACT

Petascale parallel computers with more than a million processing cores are expected to be available in a couple of years. Although MPI is the dominant programming interface today for large-scale systems that at the highest end already have close to 300,000 processors, a challenging question to both researchers and users is whether MPI will scale to processor and core counts in the millions. In this paper, we examine the issue of scalability of MPI to very large systems. We first examine the MPI specification itself and discuss areas with scalability concerns and how they can be overcome. We then investigate issues that an MPI implementation must address in order to be scalable. To illustrate the issues, we ran a number of simple experiments to measure MPI memory consumption at scale up to 131,072 processes, or 80%, of the IBM Blue Gene/P system at Argonne National Laboratory. Based on the results, we identified nonscalable aspects of the MPI implementation and found ways to tune it to reduce its memory footprint. We also briefly discuss issues in application scalability to large process counts and fea

#### - AND - ENGINEERING

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#### **Using Advanced MPI**

Modern Features of the Message-Passing Interface

William Gropp Torsten Hoefler Rajeev Thakur Ewing Lusk



# **Message Passing Interface (MPI)**

#### Message passing **libraries**:

- PVM (Parallel Virtual Machines) 1980s
- MPI (Message Passing Interface) 1990s

# **MPI = Standard API**

- Hides Software/Hardware details
- Portable, flexible
- Implemented as a library

Program				
MPI library				
Specialized Driver	Standard TCP/IP			
Custom Network HW	Standard Network HW			



### **Process Identification**

- MPI processes can be collected into groups
  - Each group can have multiple colors (some times called context)
  - Group + color == communicator (it is like a name for the group)
  - When an MPI application starts, the group of all processes is initially given a predefined name called MPI\_COMM\_WORLD
  - The same group can have many names, but simple programs do not have to worry about multiple names
- A process is identified by a unique number within each communicator, called rank
  - For two different communicators, the same process can have two different ranks: so the meaning of a "rank" is only defined when you specify the communicator



### **MPI Communicators**

 Defines the communication domain of a communication operation: set of processes that are allowed to communicate with each other.



Initially all processes are in the communicator MPI\_COMM\_WORLD.



The rank of processes are associated with (and unique within) a communicator, numbered from 0 to n-1



#### **Communicators**

mpiexec -np 16 ./test

Communicators do not need to contain all processes in the system

Every process in a communicator has an ID called as "rank"

When you start an MPI program, there is one predefined communicator MPI\_COMM\_WORLD

Can make copies of this communicator (same group of processes, but different "aliases")

The same process might have different ranks in different communicators

Communicators can be created "by hand" or using tools Simple programs typically only use the predefined communicator MPI\_COMM\_WORLD (which is sometimes considered bad practice)



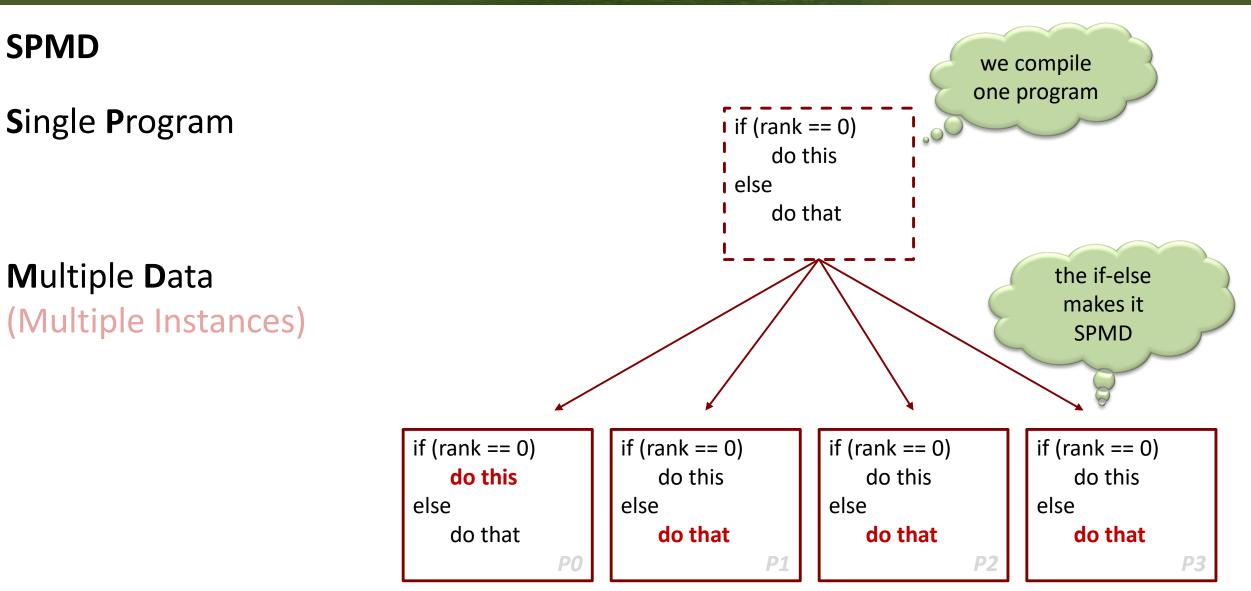
#### **Process Ranks**

Processes are identified by nonnegative integers, called *ranks* 

*p* processes are numbered *0, 1, 2, ... p-1* 

```
public static void main(String args []) throws Exception {
    MPI.Init(args);
    // Get total number of processes (p)
    int size = MPI.COMM_WORLD.Size();
    // Get rank of current process (in [0..p-1])
    int rank = MPI.COMM_WORLD.Rank();
    MPI.Finalize();
}
```





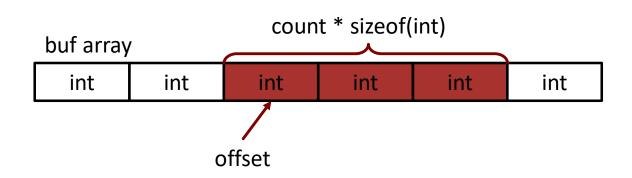
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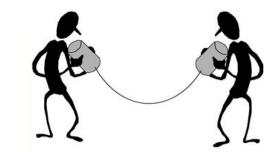


### Communication

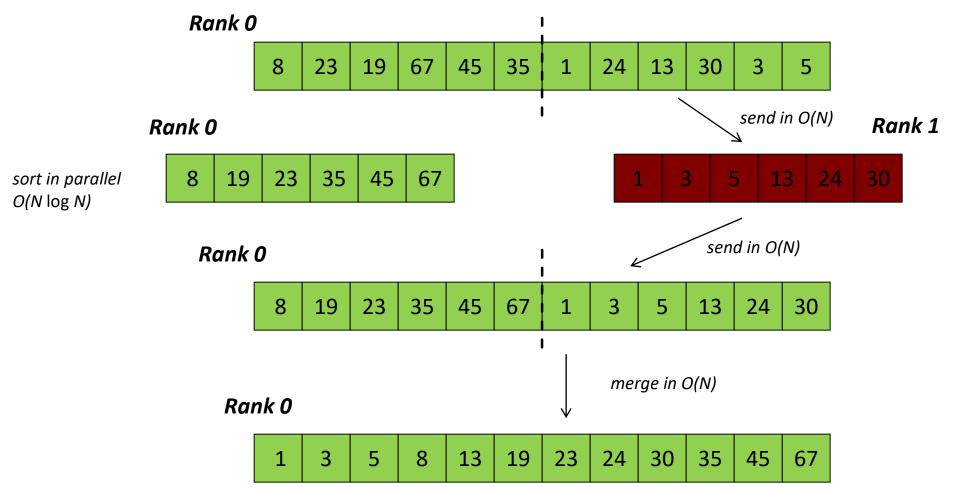
void Comm.Send(	communicator	
Object buf,	pointer to data to be sent	from MPJ Spec
int offset,		
int count,	number of items to be sent	
Datatype datatype,	data type of items, must be explicitely specified	
int dest,	destination process id	
int tag	data id tag	

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#### Parallel Sort using MPI Send/Recv



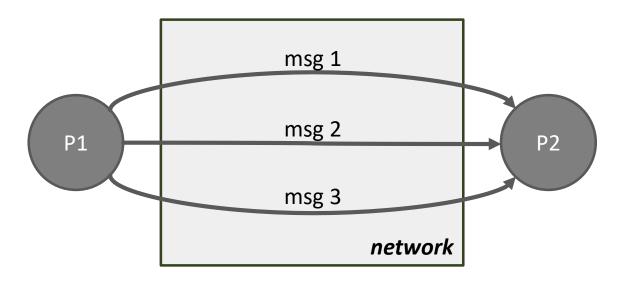
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### Message Tags

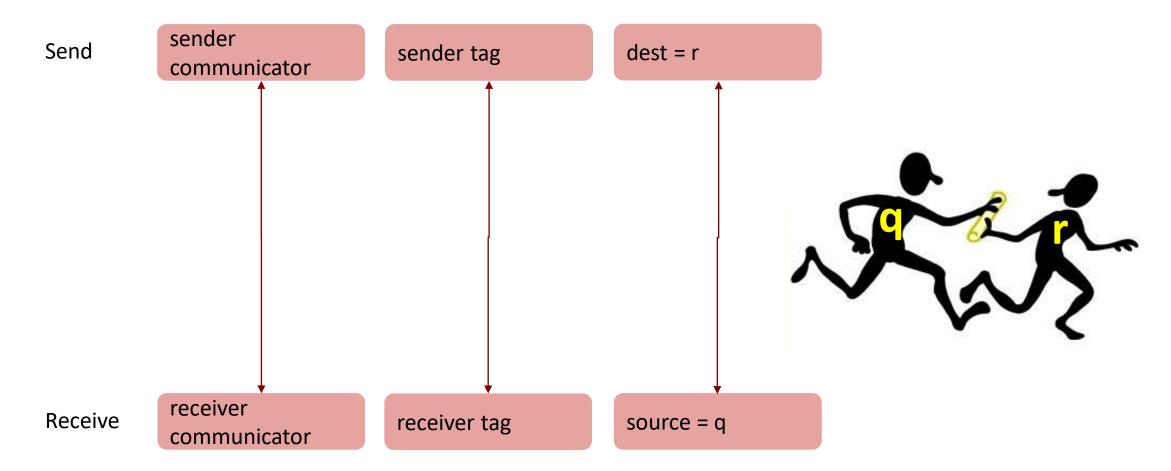
Communicating processes may need to send several messages between each other.

Message tag: differentiate between different messages being sent.





# Message matching



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### **Receiving messages**

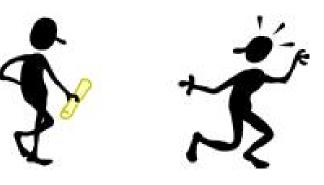
void Comm.Recv(	communicator
Object buf,	pointer to the buffer to receive to
int offset,	
int count,	number of items to be received
Datatype datatype,	data type of items, must be explicitely specified
int src,	source process id or MPI_ANY_SOURCE
int tag	data id tag or MPI_ANY_TAG
<b>\</b>	

A receiver can get a message without knowing:

- the amount of data in the message,
- the sender of the message,
- or the tag of the message.

MPI\_ANY\_SOURCE

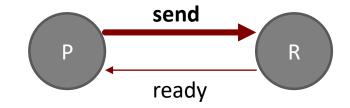
MPI\_ANY\_TAG



### **Synchronous Message Passing**

Synchronous send (Ssend)

- waits until complete message can be accepted by receiving process before completing the send
- Synchronous receive (Recv)
- waits until expected message arrives
- Synchronous routines can perform two actions
- transfer data
- synchronize processes



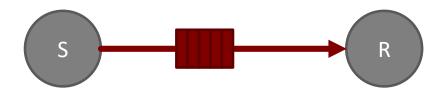


### **Asynchronous Message Passing**

Send does not wait for actions to complete before returning

• requires local storage for messages

sometimes explicit (programmer needs to care) sometimes implicit (transparent to the programmer)



In general

- no synchronisation
- allows local progress



# **Blocking / Nonblocking**

Blocking: return after *local actions* are complete, though the message transfer may not have been completed

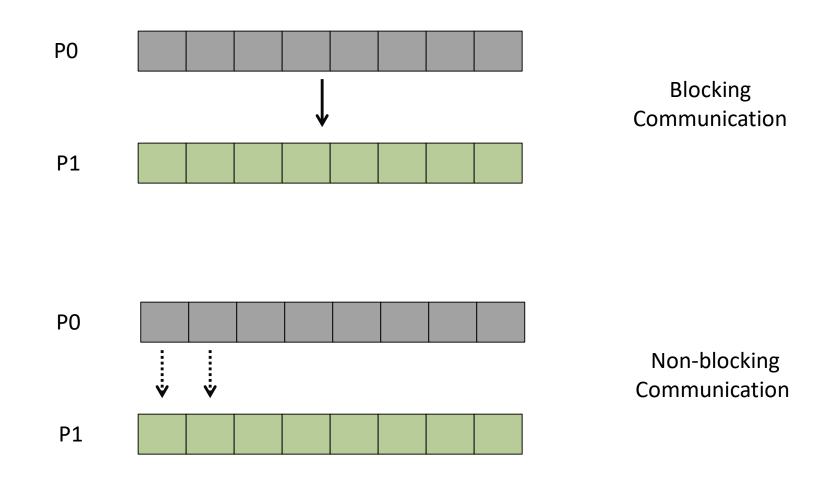
Non-blocking: return immediately

- sometimes assumes that data storage to be used for transfer is not modified by subsequent statements until transfer complete
- sometimes implementation dependent local buffers are or have to be provided





### A Non-Blocking communication example



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# Synchronous / Asynchronous vs Blocking / Nonblocking

# Synchronous / Asynchronous

about communication between sender and receiver

# **Blocking / Nonblocking**

about local handling of data to be sent / received



# **MPI Send and Receive Defaults**

# Send

- blocking,
- synchrony implementation dependent

Danger of Deadlocks. Don't make any assumptions!

depends on existence of buffering, performance considerations etc

### Recv

blocking

There are a lot of different variations of this in MPI.





# **Sources of Deadlocks**

- Send a large message from process 0 to process 1
  - If there is insufficient storage at the destination, the send must wait for the user to provide the memory space (through a receive)
- What happens with this code?

Process 0	Process 1	
Send(1)	Send(0)	
Recv(1)	Recv(0)	

• This is called "unsafe" because it depends on the availability of system buffers in which to store the data sent until it can be received



### Some Solutions to the "unsafe" Problem

Order the operations more carefully:

Process 0	Process 1
Send(1)	Recv(0)
Recv(1)	Send(0)

• Supply receive buffer at same time as send:

Process 0	Process 1
Sendrecv(1)	Sendrecv(0)



### More Solutions to the "unsafe" Problem

Supply own space as buffer for send

	Process 0	Process 1	
	Bsend(1) Recv(1)	Bsend(0) Recv(0)	
Use	non-blocking	operations:	
	Process 0	Process 1	
	Isend(1) Irecv(1) Waitall	Isend(0) Irecv(0) Waitall	



# **MPI is Simple**

Many parallel programs can be written using just these six functions, only two of which are non-trivial:

The second second

- MPI\_INIT initialize the MPI library (must be the first routine called)
- MPI\_COMM\_SIZE get the size of a communicator
- MPI\_COMM\_RANK get the rank of the calling process in the communicator
- MPI\_SEND send a message to another process
- MPI\_RECV send a message to another process
- MPI\_FINALIZE clean up all MPI state (must be the last MPI function called by a process)
- For performance, however, you need to use other MPI features

# **Example: compute Pi**

- The irrational number Pi has many digits
  - And it's not clear if they're randomly distributed!
- But they can be computed

$$\pi \approx h \sum_{i=0}^{N-1} \frac{4}{1 + (h(i + \frac{1}{2}))^2}$$

for(int i=0; i<numSteps; i++) {
 double x=(i + 0.5) \* h;
 sum += 4.0/(1.0 + x\*x);
}
double pi=h \* sum ;</pre>

### Pi record smashed as team finds twoquadrillionth digit

By Jason Palmer Science and technology reporter, BBC News

() 16 September 2010 | Technology

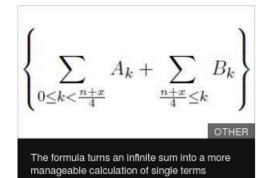
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A researcher has calculated the 2,000,000,000,000,000,000 digit of the mathematical constant pi - and a few digits either side of it.

Nicholas Sze, of tech firm Yahoo, said that when pi is expressed in binary, the two quadrillionth "bit" is 0.

Mr Sze used Yahoo's Hadoop cloud computing technology to more than double the previous record.

It took 23 days on 1,000 of Yahoo's computers - on a standard PC, the calculation would have taken 500 years.





# **Pi's parallel version**

```
MPI.Init(args);
... // declare and initialize variables (sum=0 etc.)
int size = MPI.COMM WORLD.Size();
int rank = MPI.COMM WORLD.Rank();
for(int i=rank; i<numSteps; i=i+size) {</pre>
   double x=(i + 0.5) * h;
   sum += 4.0/(1.0 + x^*x);
}
if (rank != 0) {
   double [] sendBuf = new double []{sum};
   // 1-element array containing sum
   MPI.COMM WORLD.Send(sendBuf, 0, 1, MPI.DOUBLE, 0, 10);
}
else { // rank == 0
   double [] recvBuf = new double [1] ;
   for (int src=1 ; src<P; src++) {</pre>
      MPI.COMM_WORLD.Recv(recvBuf, 0, 1, MPI.DOUBLE, src, 10);
      sum += recvBuf[0];
double pi = h * sum; // output pi at rank 0 only!
MPI.Finalize();
```

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# **COLLECTIVE COMMUNICATION**



### **Group Communication**

Up to here: point-to-point communication

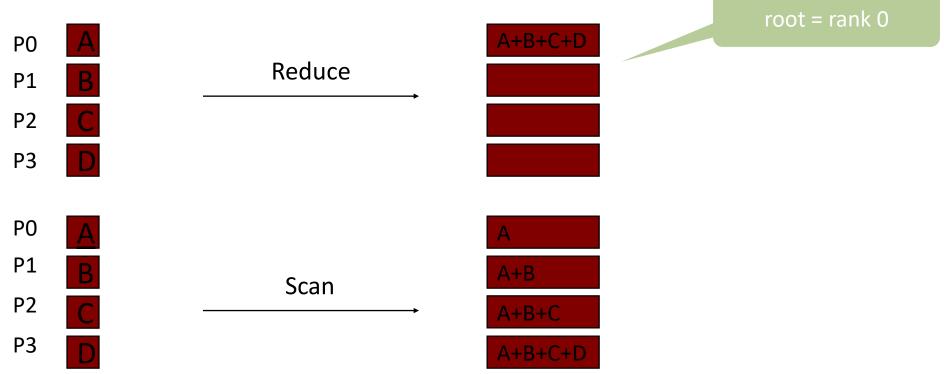
MPI also supports communications among groups of processors

- not absolutely necessary for programming (but very nice!)
- but essential for performance

Examples: broadcast, gather, scatter, reduce, barrier, ...

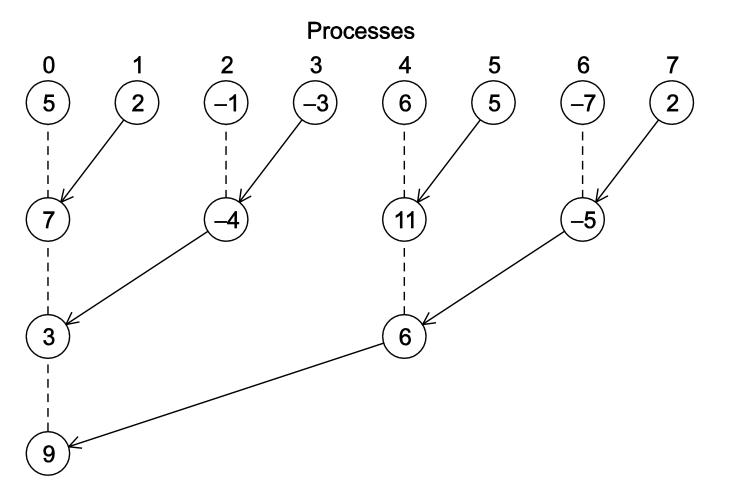
### **Collective Computation - Reduce**

#### 



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### **Reduce implementation: a tree-structured global sum**



- In the first phase:

   (a) Process 1 sends to 0, 3 sends to
   2, 5 sends to 4, and 7 sends to 6.
   (b) Processes 0, 2, 4, and 6 add in the received values.
- 2. Second phase:
  (c) Processes 2 and 6 send their new values to processes 0 and 4, respectively.
  (d) Processes 0 and 4 add the received values into their new values.
- 3. Finally:

(a) Process 4 sends its newest value to process 0.

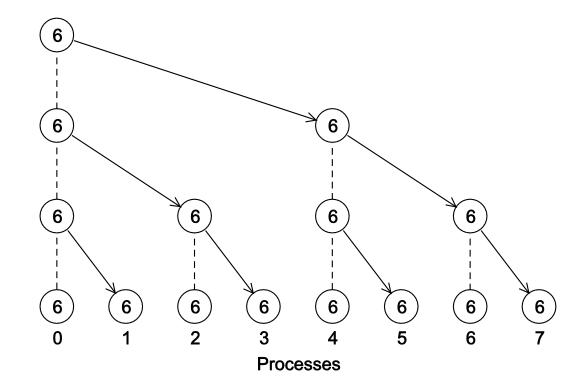
(b) Process 0 adds the received value to its newest value.



### **Collective Data Movement - Broadcast**

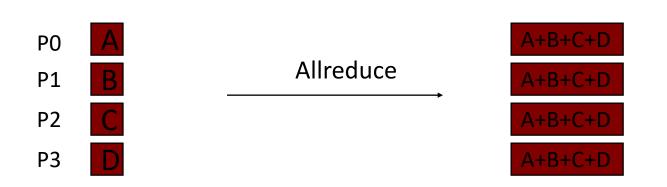


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### **Collective Computation - Allreduce**

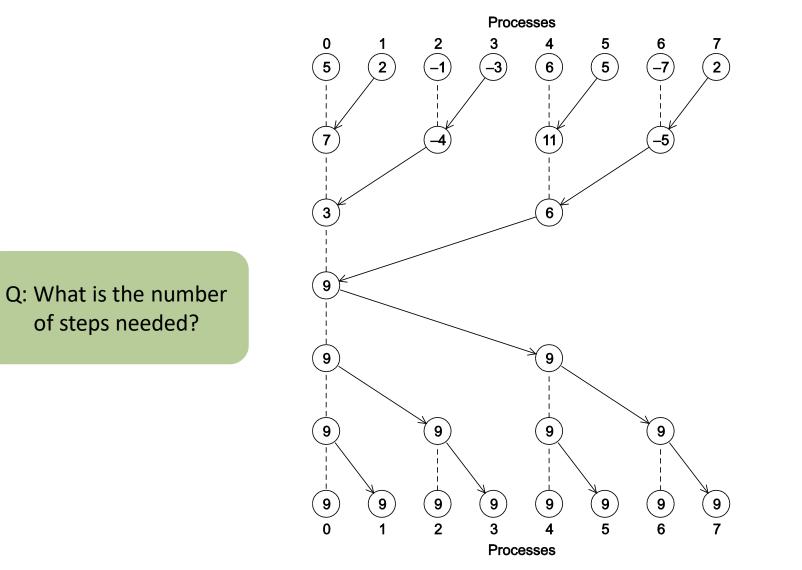
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Useful in a situation in which all of the processes need the result of a global sum in order to complete some larger computation.



### Allreduce = Reduce + Broadcast?

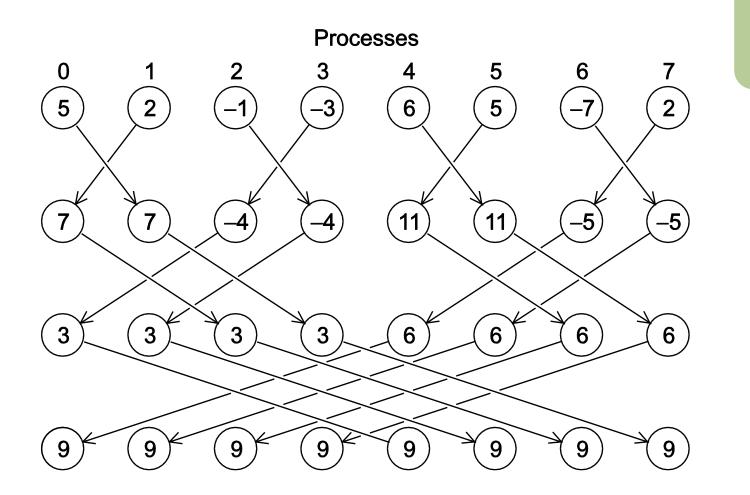


- Constanting

A global sum followed by distribution of the result.



### Allreduce ≠ Reduce + Broadcast



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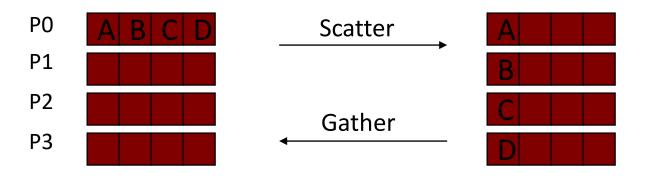
Q: What is the number of steps needed?



A butterfly-structured global sum.



# **Collective Data Movement – Scatter/Gather**

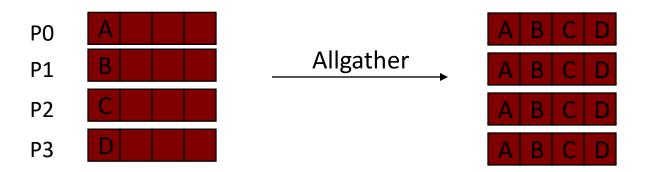


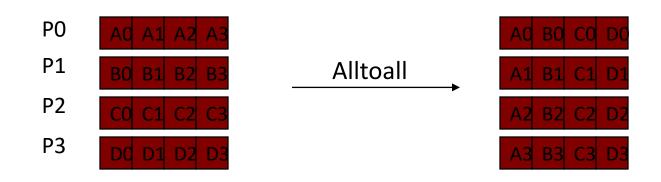
- Scatter can be used in a function that reads in an entire vector on process 0 but only sends the needed components to each of the other processes.
- Gather collects all of the components of the vector onto destination process, then destination process can process all of the components.



### More Collective Data Movement – some more (16 functions total!)

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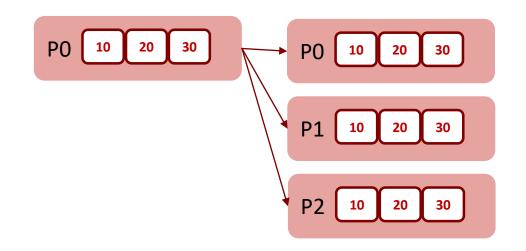




**Compute** 
$$y = A \cdot x$$
, *e.g.*,  $A = \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$   $x = \begin{bmatrix} 10 \\ 20 \\ 30 \end{bmatrix}$   $y = \begin{bmatrix} A_{1.} \cdot x \\ A_{2.} \cdot x \\ A_{3.} \cdot x \end{bmatrix}$ 

States and

### 1. Broadcast x

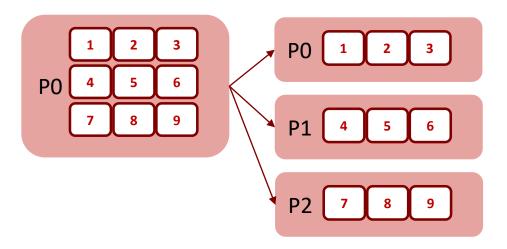




**Compute** 
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### 2. Scatter A

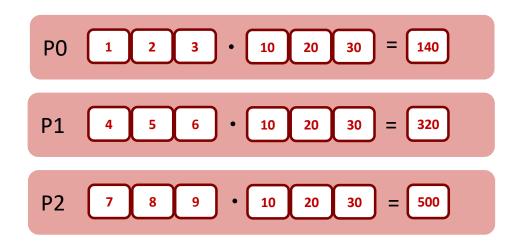




**Compute** 
$$y = A \cdot x$$
, *e.g.*  $A = \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$   $x = \begin{bmatrix} 10 \\ 20 \\ 30 \end{bmatrix}$   $y = \begin{bmatrix} A_{1} \cdot x \\ A_{2} \cdot x \\ A_{3} \cdot x \end{bmatrix}$ 

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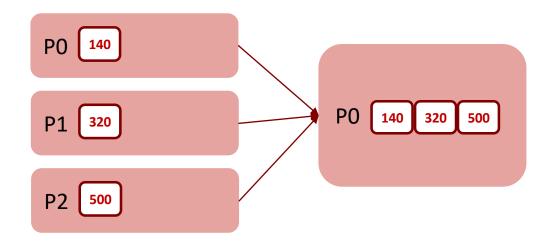
### **3. Compute locally**





**Compute** 
$$y = A \cdot x$$
, *e.g.*  $A = \begin{bmatrix} 1 & 2 & 3 \\ 4 & 5 & 6 \\ 7 & 8 & 9 \end{bmatrix}$   $x = \begin{bmatrix} 10 \\ 20 \\ 30 \end{bmatrix}$   $y = \begin{bmatrix} A_{1.} \cdot x \\ A_{2.} \cdot x \\ A_{3.} \cdot x \end{bmatrix}$ 

### 4. Gather result y





### Iterations

Assume we want to apply the matrix-vector product iteratively

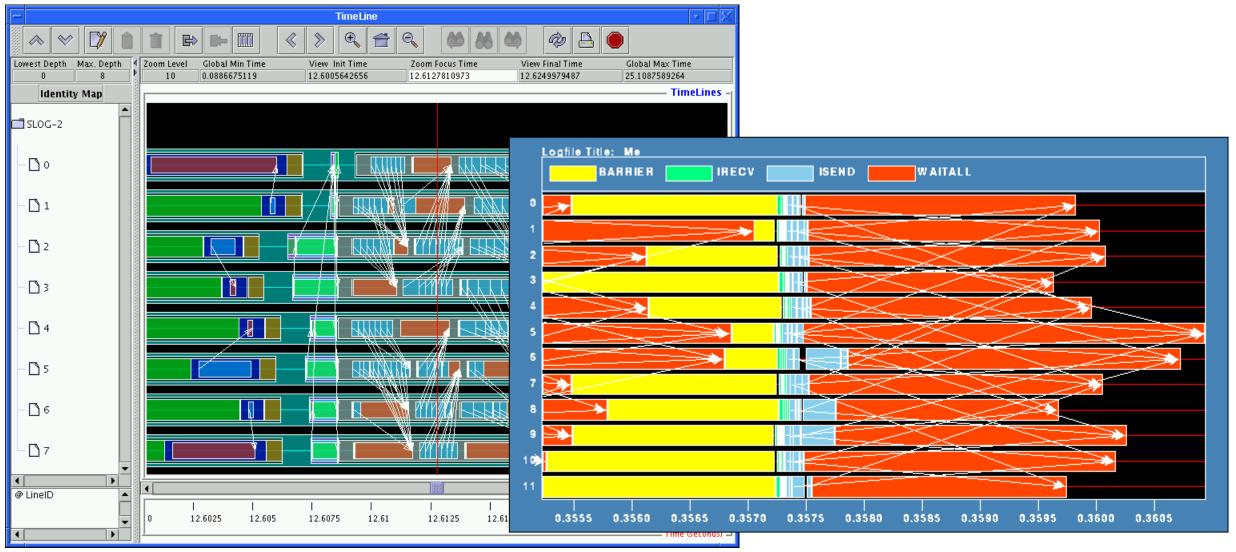
$$y_n = A y_{n-1}$$

Example Application:

Eigenvalue Problem for Probability Matrix, as used in Google's Pagerank algorithm.

Then each process needs the results of other processes after one step.  $\rightarrow$  Need for Gather + Broadcast in one go.

### **Visualizing Program Behavior**



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# **MPI conclusion**

- The de-facto interface for distributed parallel computing (nearly 100% market share in HPC)
- Elegant and simple interface
  - Definitely simpler than shared memory (no races, limited conflicts, avoid deadlocks with nonblocking communication)
- We only covered the basics here, MPI-3.1 (2015) has 600+ functions
  - More concepts:
    - Derived Datatypes
    - Process Topologies
    - Nonblocking and neighborhood collectives

One-sided accesses (getting the fun of shared memory back ...) Profiling interfaces