

# ME 5311: Week 14

Today

## **Introduction to parallel computing**

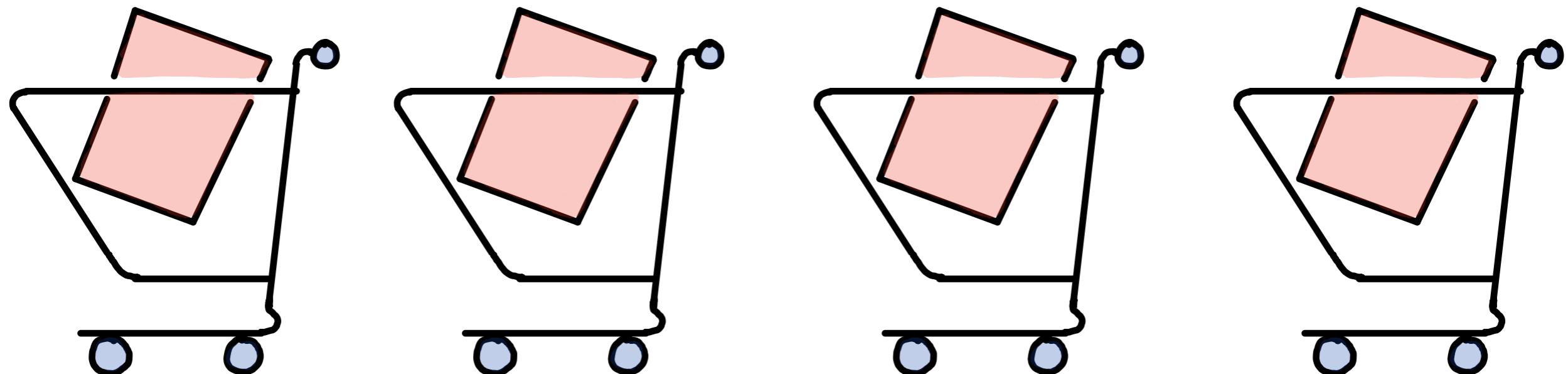
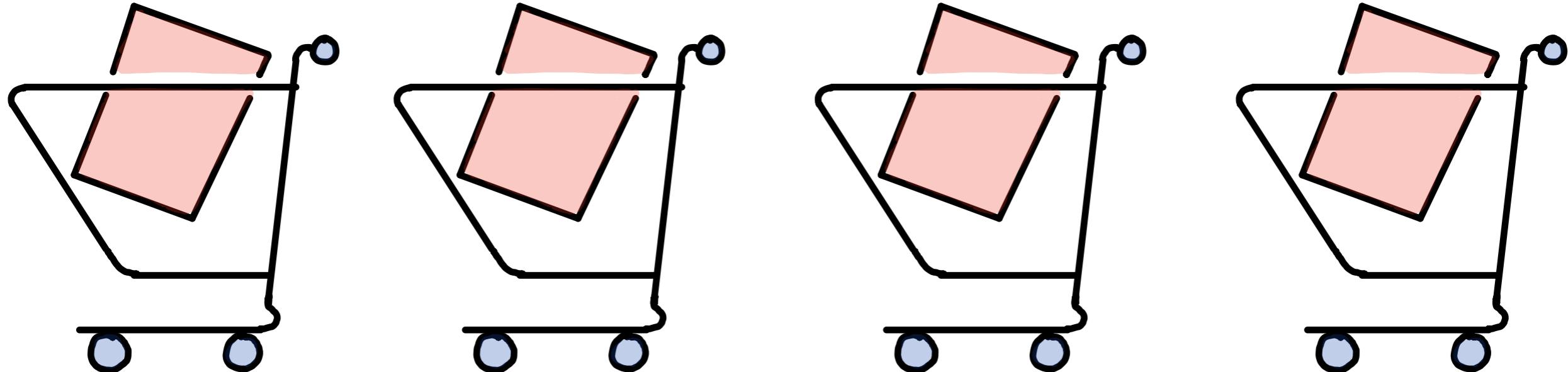
- Goal: Understand basic principles of parallel computing for CFD
- Introduce and discuss the two most common approaches
- Get a flavor of “real” programming

# What is parallel computing?

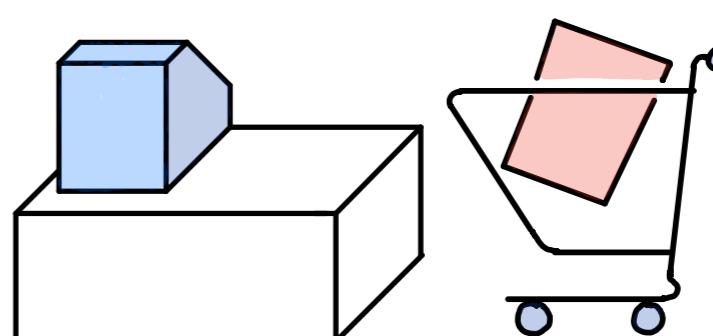
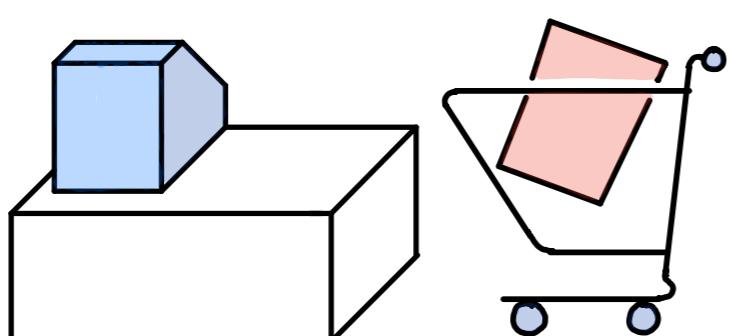
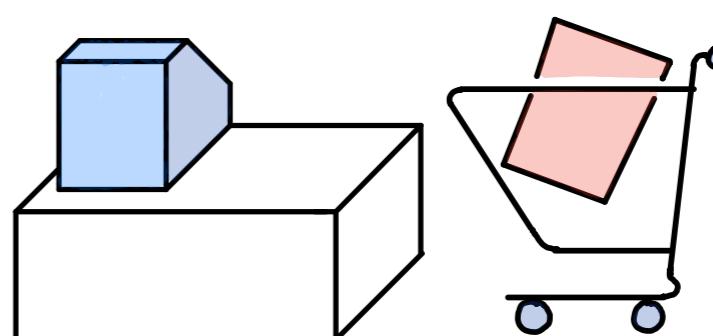
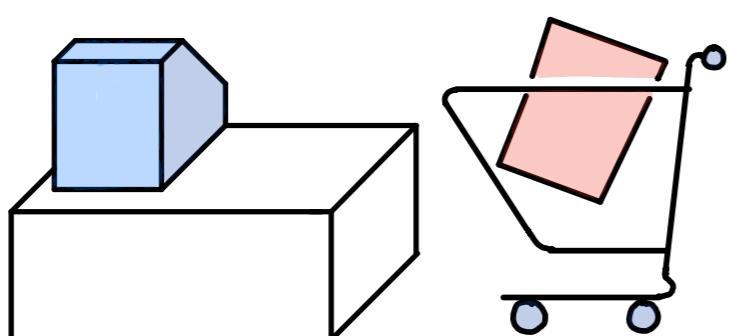
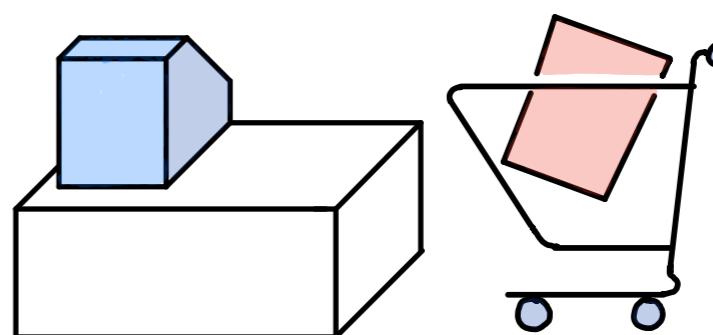
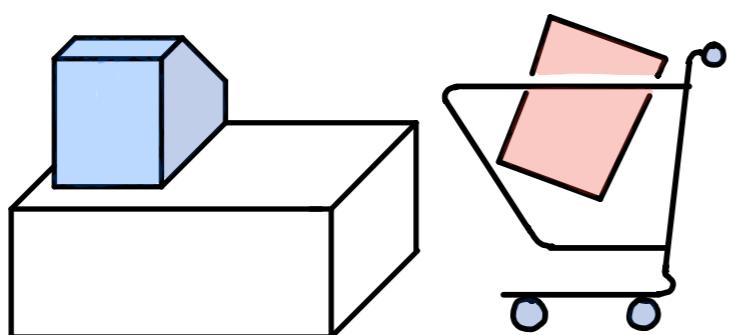
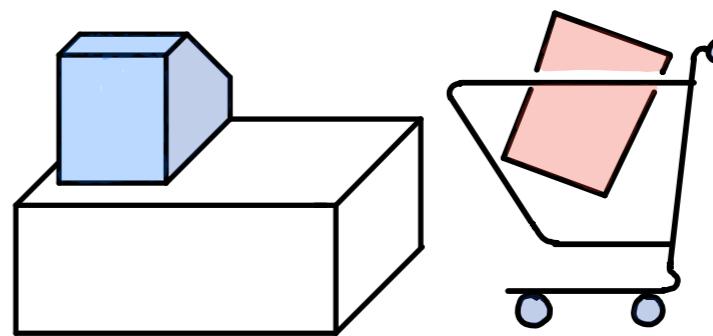
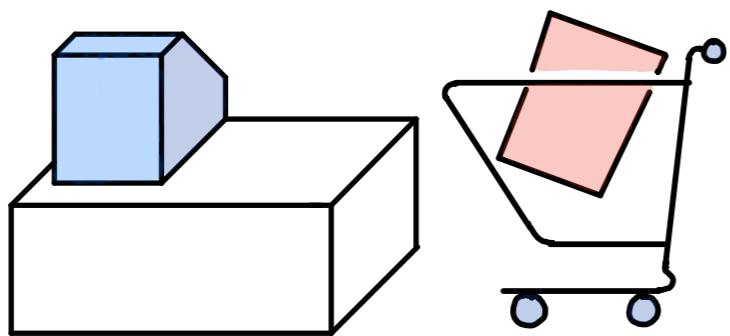


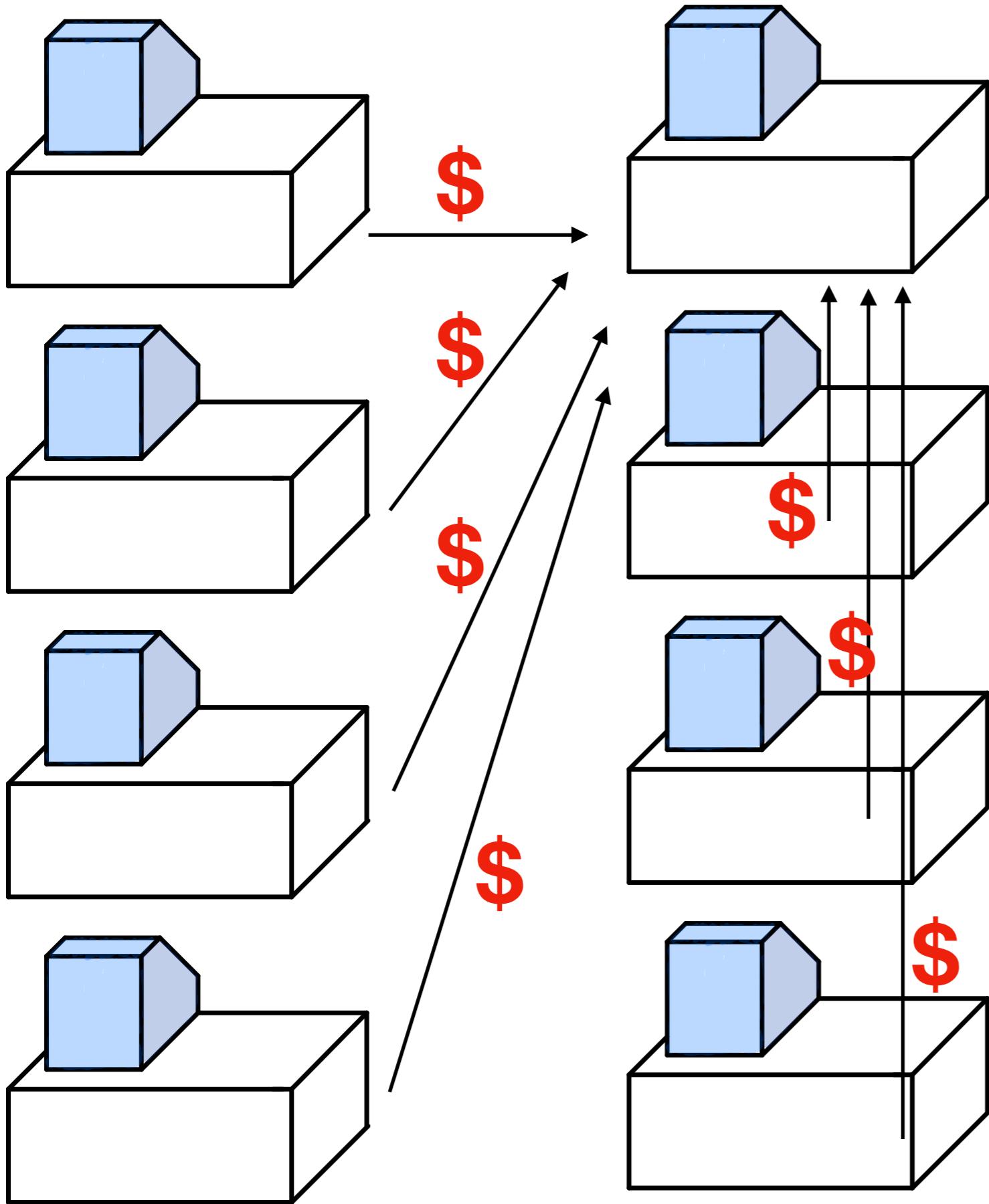
# Need for parallel computing

- We would like to solve a “big” problem
- “Big” can mean:
  - Takes a long time to compute: split the work on multiple CPUs
  - Requires a lot of computer memory: use memory from different “computer nodes”

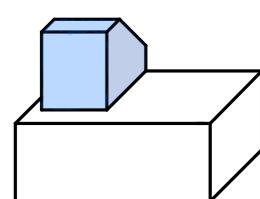
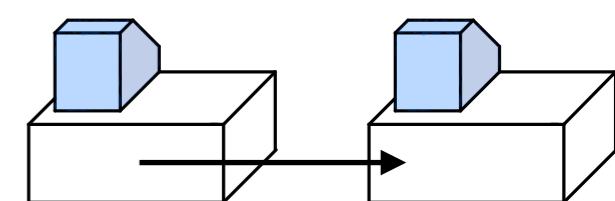
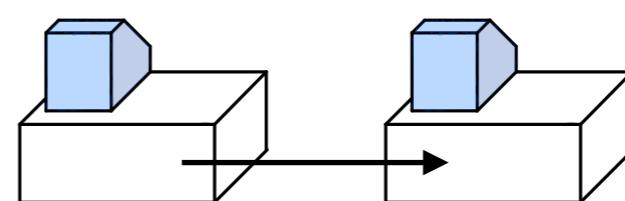
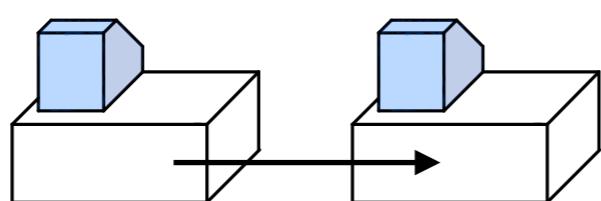
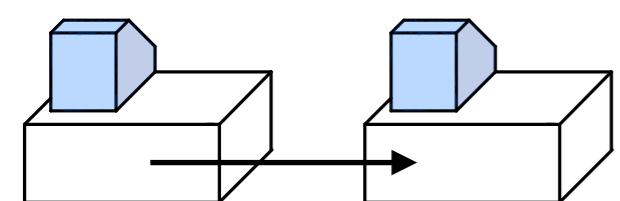


# Splitting the work: granularity





**Processes must:**  
**communicate**  
**coordinate**  
**combine results**

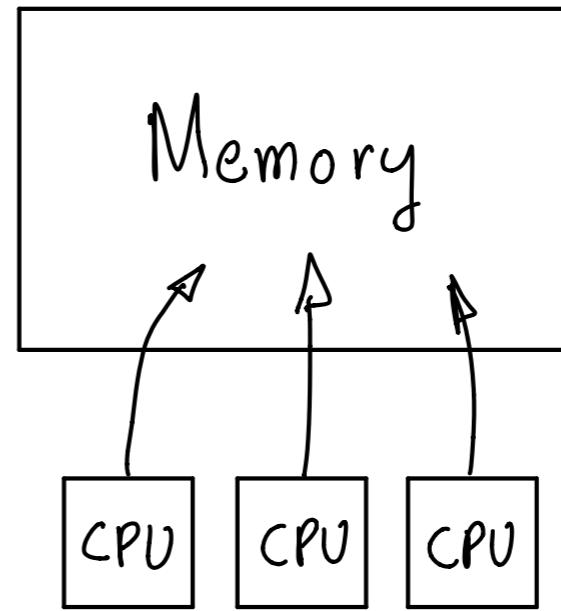


# Basic principle: Program copies itself

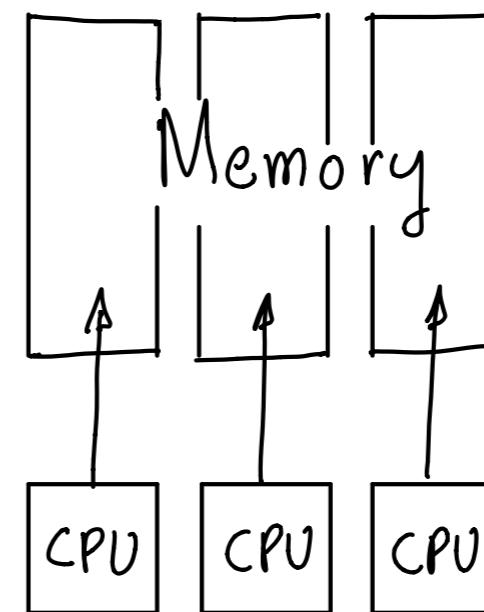


# Traditional parallel models

- **Shared memory:**  
OpenMP, Threads

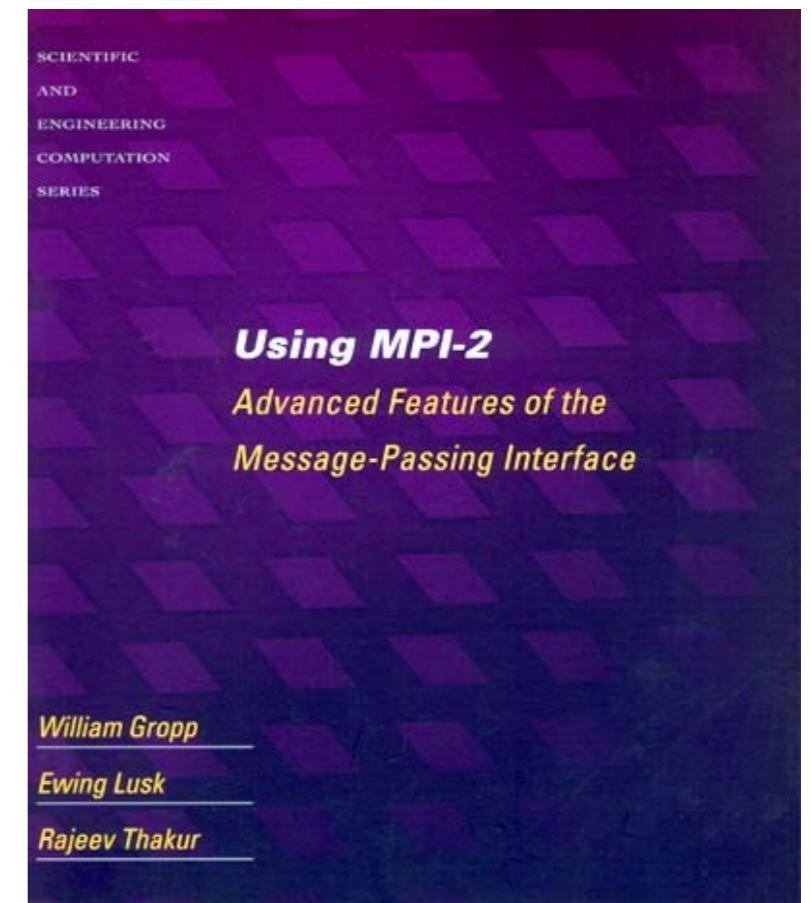
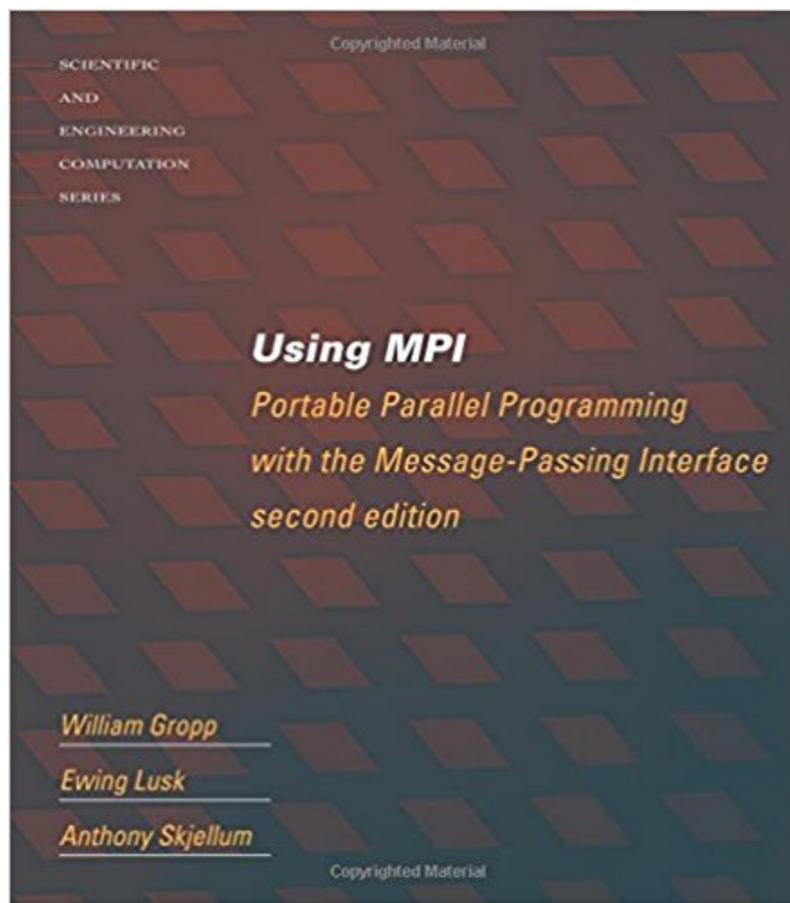


- **Described memory:**  
Message Passing Interface



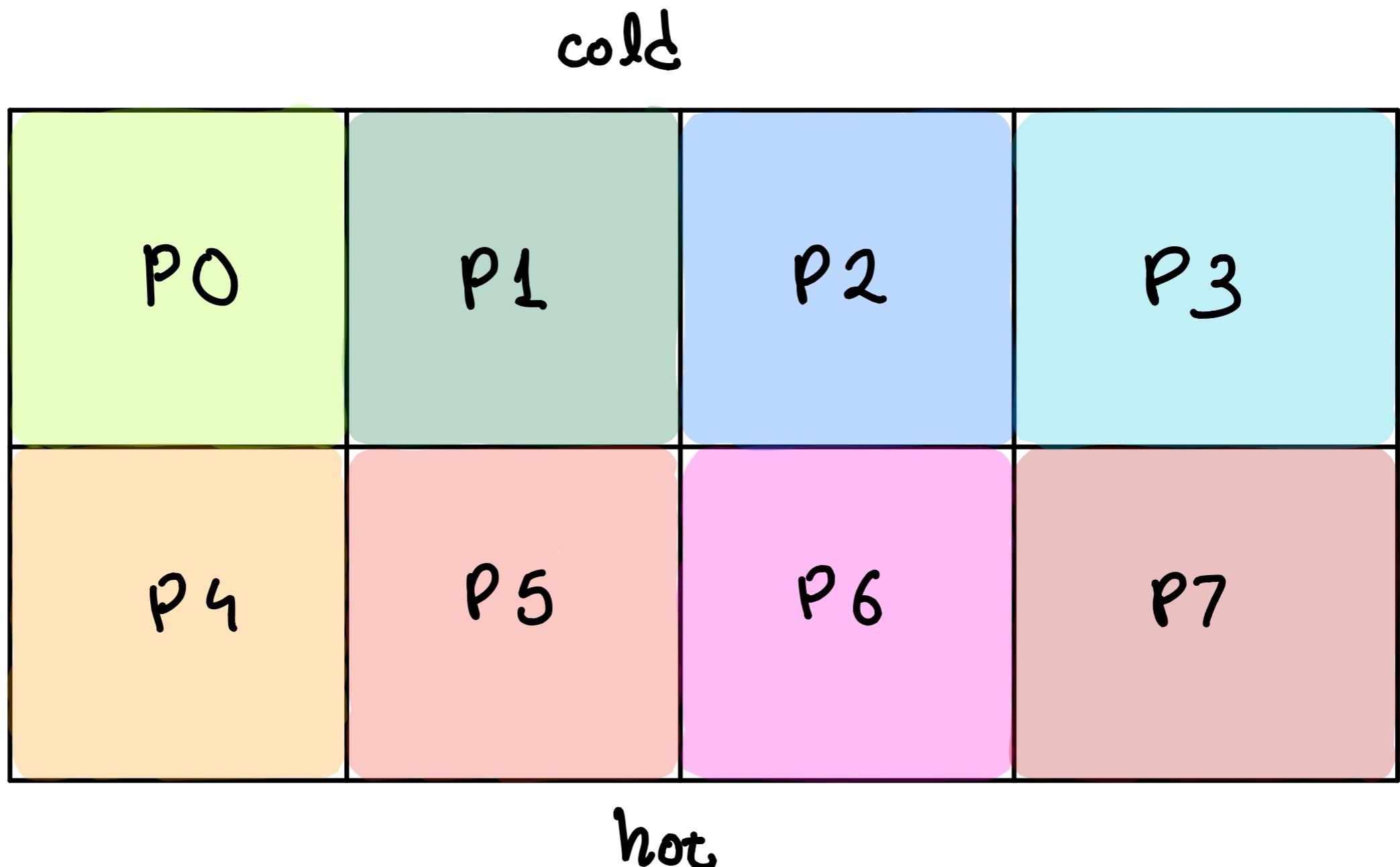
# Implementations

**Goal is to “tell” the computer what you want to do...  
e.g., send data from X to Z**

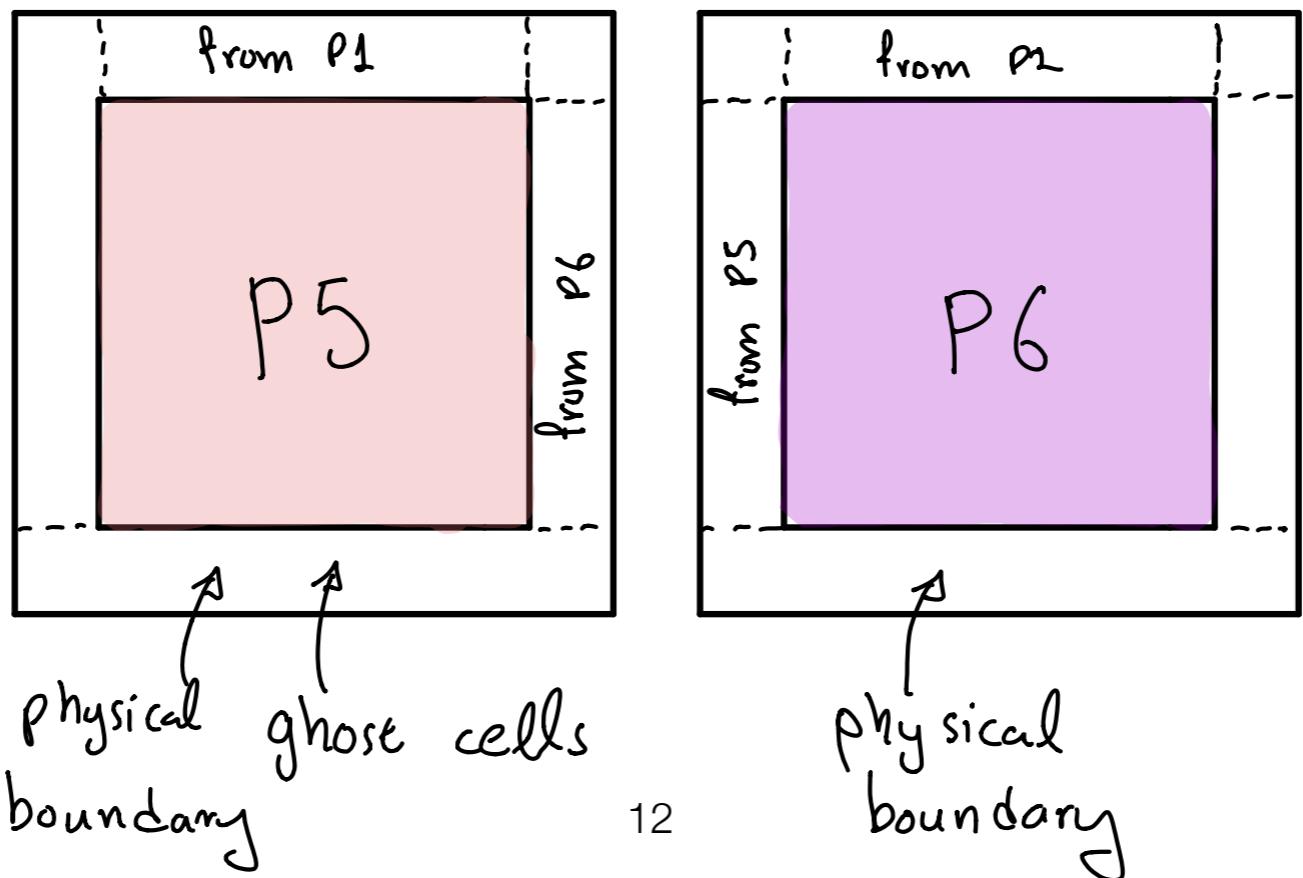
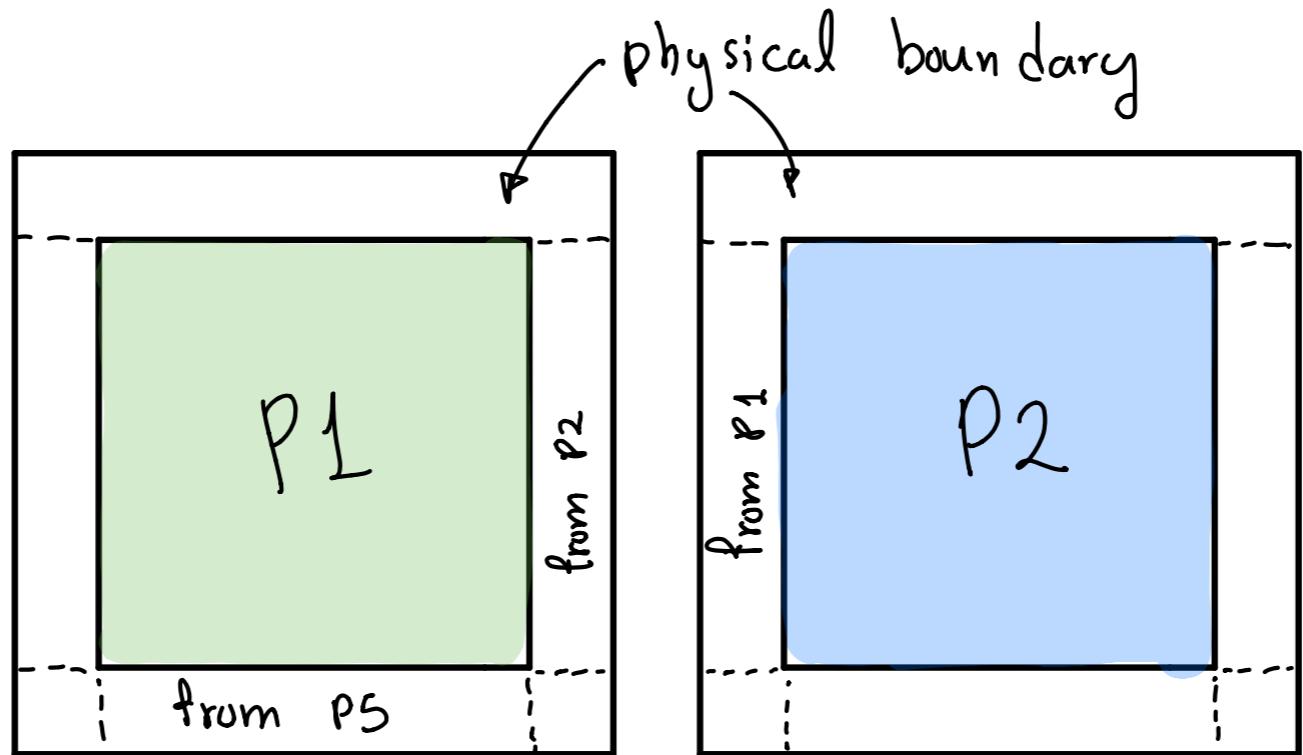


**Implementation literature/books assume's reader is experienced programmer**

# Computational domain decomposition



# Process communication



# Non-blocking communication

```
call mpi_isend(var(1,n2-2*nghost(1)+1,nghost(2)+1),1,xstride,  
pxfwd, 130, MPI_COMM_WORLD, req(1), ierror)  
  
call mpi_irecv(var(1,1,nghost(2)+1), 1, xstride, pxback, 130,  
MPI_COMM_WORLD, req(4), ierror)  
  
call mpi_isend(var(1,nghost(1)+1,nghost(2)+1), 1, xstride, pxback,  
140, MPI_COMM_WORLD, req(2), ierror)  
  
call mpi_irecv(var(1,n2-nghost(1)+1,nghost(2)+1), 1, xstride,  
pxfwd, 140, MPI_COMM_WORLD, req(3), ierror)  
  
...  
  
call mpi_waitall(16,req,stats,ierror)
```