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# Parallel Computing

## WS 2017/18

### Session 6: MPI Recap, Stencil Operations

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# Two-sided vs. One-sided

## Two-sided

- **Memory is private** to each process.
- When the sender calls the MPI\_Send operation and the receiver calls the MPI\_Recv operation, data in the sender memory is copied to a buffer then sent over the network, where it is copied to the receiver memory.
- **Drawback:** sender has to wait for the receiver to be ready to receive the data before it can send the data.
- Both sender and receiver have to state a specific call for the communication.

→ **Coupled program flow**

## One-sided

- Sections in local memory are made **accessible among processes**.
- Requires **only one process to transfer data**, decouples data transfer from system synchronization.
- **MPI 3.0** supports one-sided passive target communication **without the intervention of the remote process** via Remote Direct Memory Access (RDMA).

That is: send or receive data without any local action.

→ **Decoupled program flow**

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→ **Decoupled program flow**

In the real world, passive RDMA in most MPI implementations is either buggy, or inefficient (barrier spin-locks and other delightful hacks) or both, but it's getting better.

## One-sided Operations

### Standard

MPI\_Put

MPI\_Get

MPI\_Accumulate

### Request-based

MPI\_Rput

MPI\_Rget

MPI\_Raccumulate

- All data movement operations are **non-blocking**.
- **Requires explicit synchronization call** to ensure completion, e.g.:

MPI\_Wait(req)

MPI\_Win\_fence(win)

MPI\_Win\_flush(win)

...

## MPI\_Get / MPI\_Put

Origin: calling (i.e. local) process

Target: remote process

```
MPI_Get(oaddr, ocount, otype,  
        trank, tdisp, tcount, ttype, window)
```

Transfer elements from target in window[tdisp:tcount]  
into local buffer oaddr at origin.

```
MPI_Put(oaddr, ocount, otype,  
        trank, tdisp, tcount, ttype, window)
```

Transfer elements from local buffer oaddr at origin  
to target into window[tdisp:tcount].

## (Blocking / Nonblocking) x (One-sided x Two-sided)

	Blocking	Nonblocking
Two-sided	MPI_Send	MPI_Isend
One-sided	MPI_Put	MPI_Rput

One-sided communication can be used to implement collective operations.

Pop quiz:           How would you implement a reduce operation using one-sided communication?

Example:            **Find minimum value in distributed array.**



## One-sided true passive RMA Example

```
MPI_Comm comm = MPI_COMM_WORLD;
// Create local window buffer of 4096 integers:
int * winbuf;
MPI_Alloc_mem(sizeof(int)*4096, MPI_INFO_NULL, &winbuf);
// Create window, collective operation:
MPI_Win win;
MPI_Win_create(winbuf, sizeof(int)*4096, sizeof(int),
               MPI_WIN_INFO_NULL, comm, &win);
// Start passive RMA epoch, collective operation:
MPI_Win_lock_all(0, win);

...
```

## One-sided true passive RMA Example ctd.

```
// rank_0.buf[0:99] <- rank_1.win[1024:1123]
```

rank 1:

```
-----+-----  
for (int i=0; i<128; i++)  
    winbuf[i+1024] = 42;  
}  
MPI_Win_flush_all(win);
```

rank 0:

```
-----+-----  
int buf[100];  
MPI_Get(buf, 100, MPI_INT,  
        1, 1024, comm);  
MPI_Win_flush(1, win);
```

**What could go wrong?**

## One-sided true passive RMA Example ctd.

```
// rank_0.buf[0:99] <- rank_1.win[1024:1123]
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rank 1:

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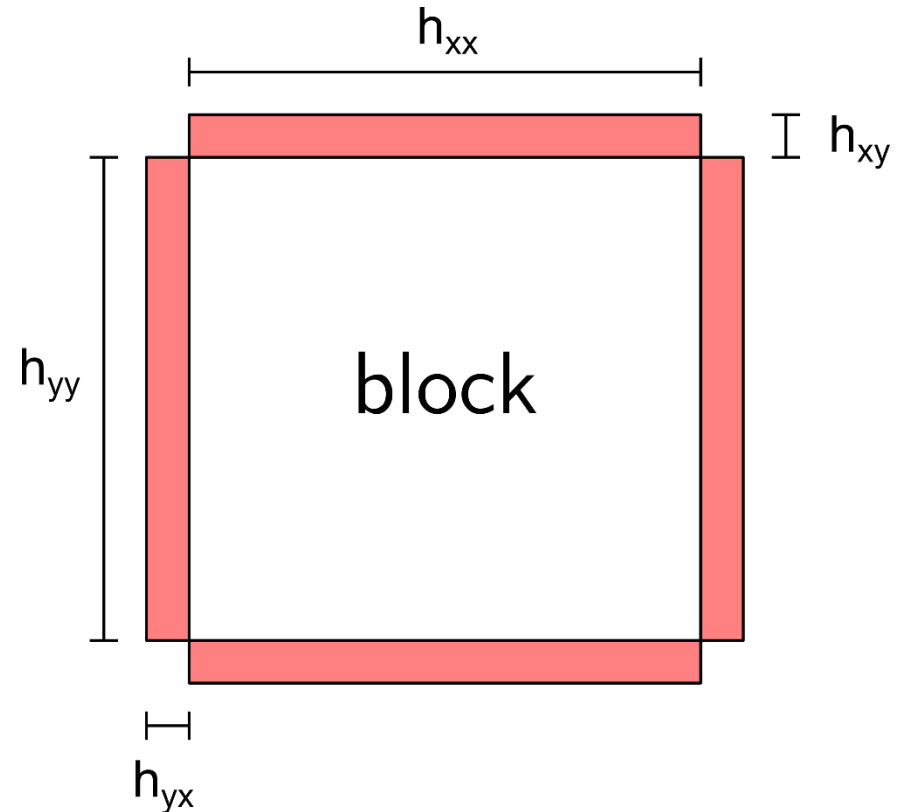
**Data races, as known from multi-threading.**

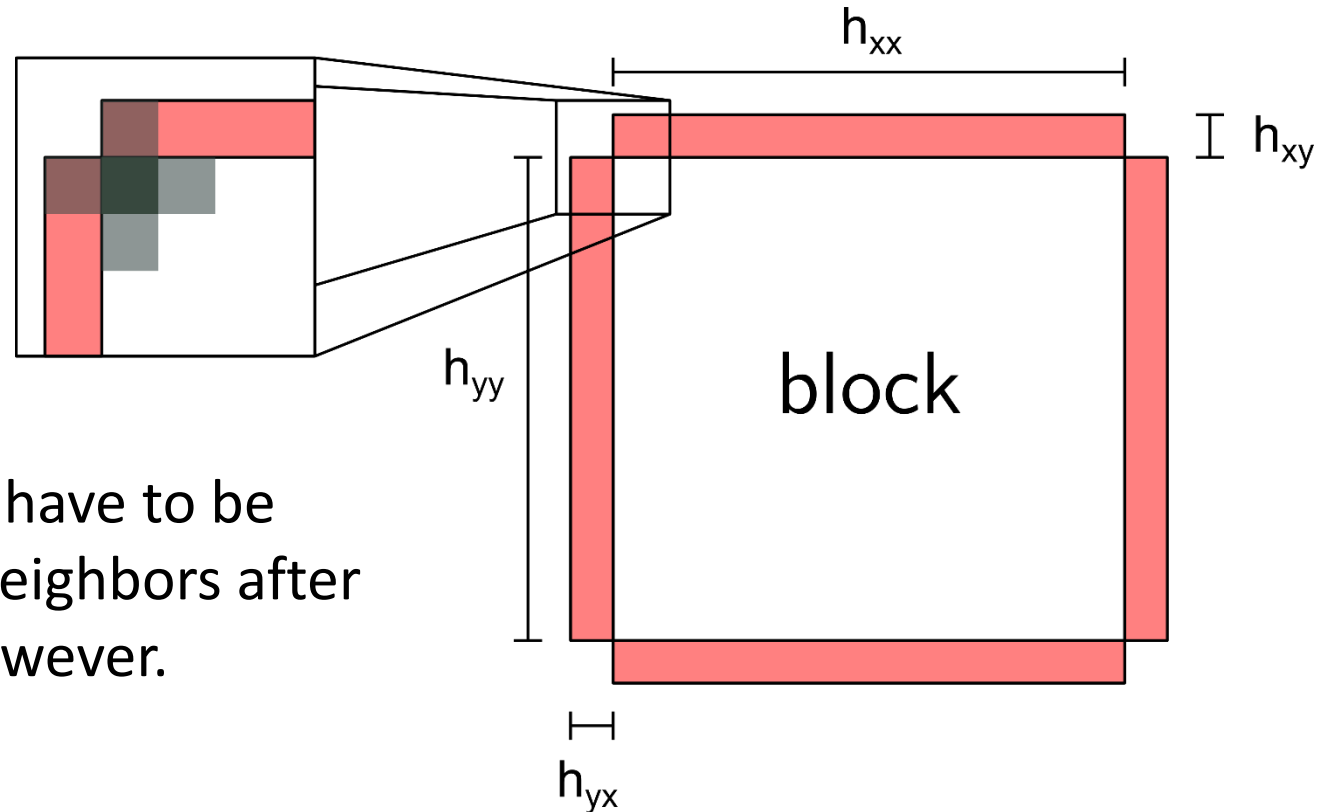
# Stencil Operations



rank 0 block [0,0]	rank 1 block [0,1]
rank 2 block [1,0]	rank 3 block [1,1]

Computing the inner-most values in a local block is straight-forward (local-only).





Ghost cells (halo) have to be exchanged with neighbors after each iteration, however.

Assuming a 1-nn stencil,

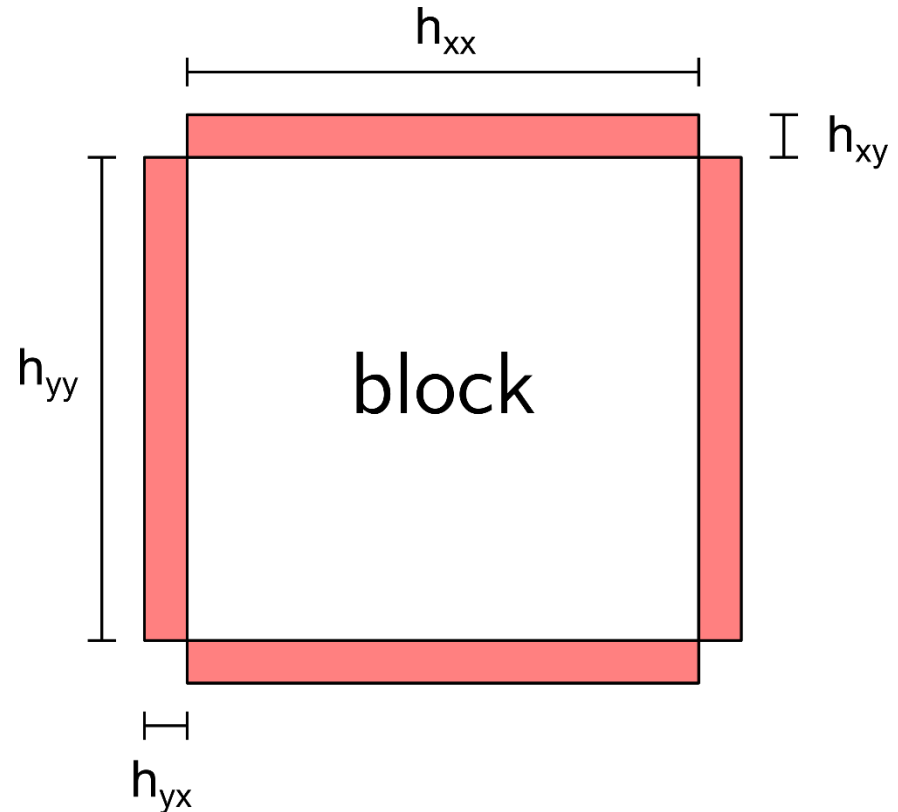
block size  $B = b_x \times b_y$

field size  $N = n_x \times n_y$

Elements exchanged with all  
neighbors per block:

$4b_x b_y$

**Surface-to-volume ratio?**





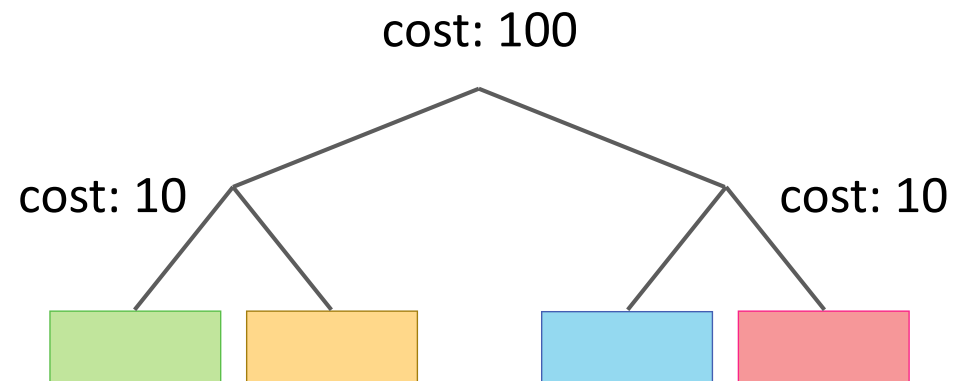
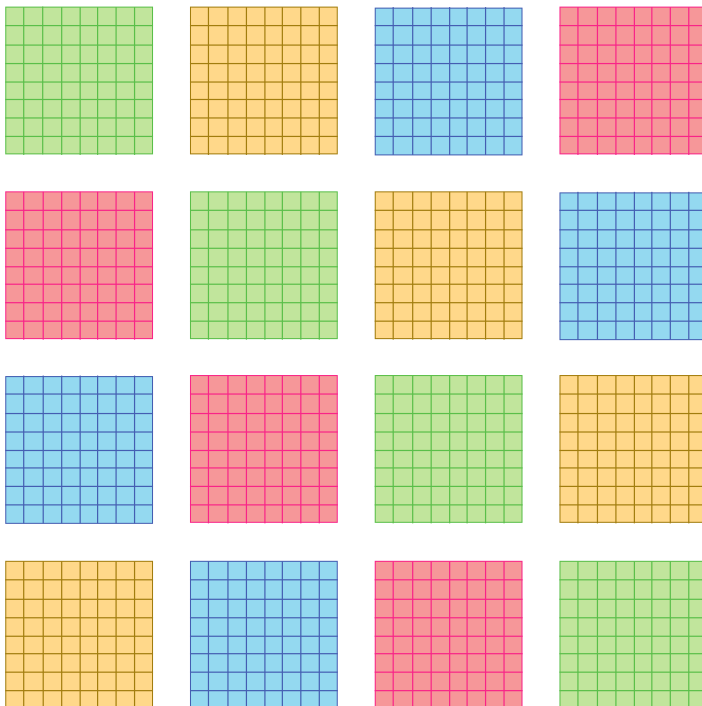
## Surface/Volume Ratio

- For high degree of parallelism: **select small block size** (more processes → more blocks → smaller block size)
- But: small block size affects border exchanges, surface/volume ratio increases.
- As the size of a block increases its **volume grows faster than its surface area**.  
Square-Cube Law:  $O(n^3)$  vs  $O(n^2)$
- High ratio → more the time spent on communication per iteration, less time left to spend on actual computations.

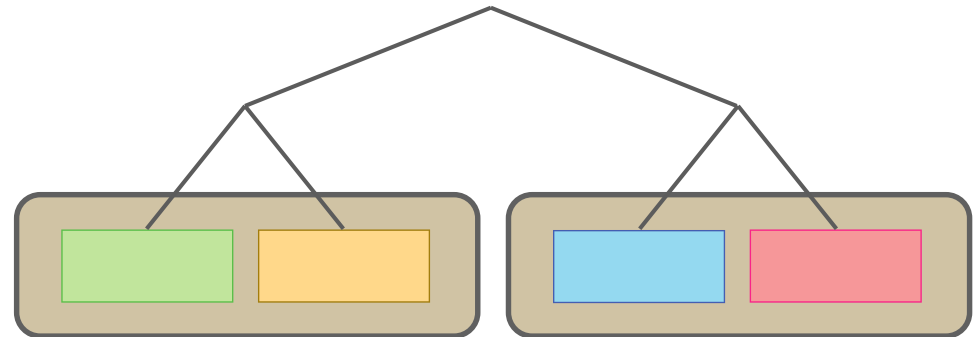
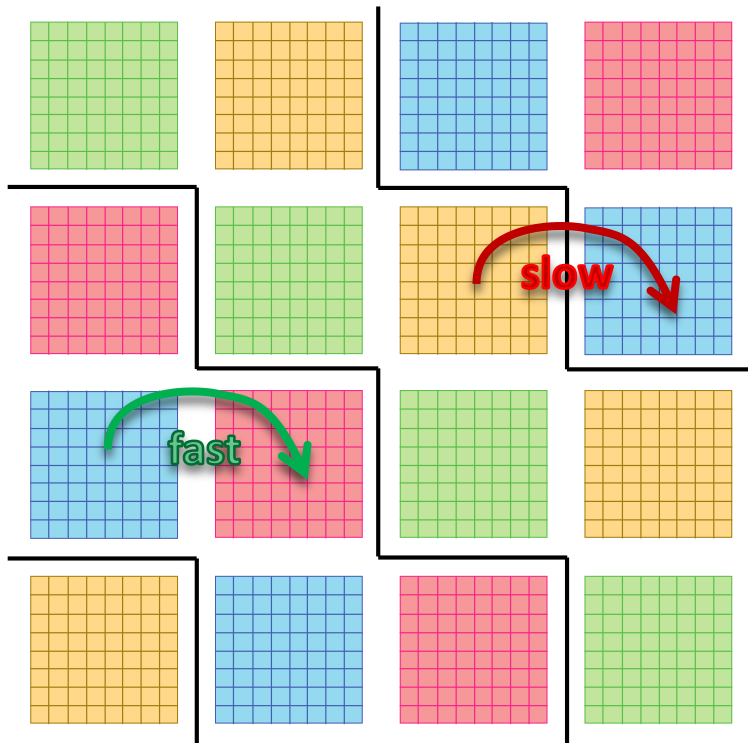
# **Domain Decomposition vs. Process Topology vs. Hardware Locality**



Data distribution, hardware and communication patterns are  
**highly interdependent with respect to performance**

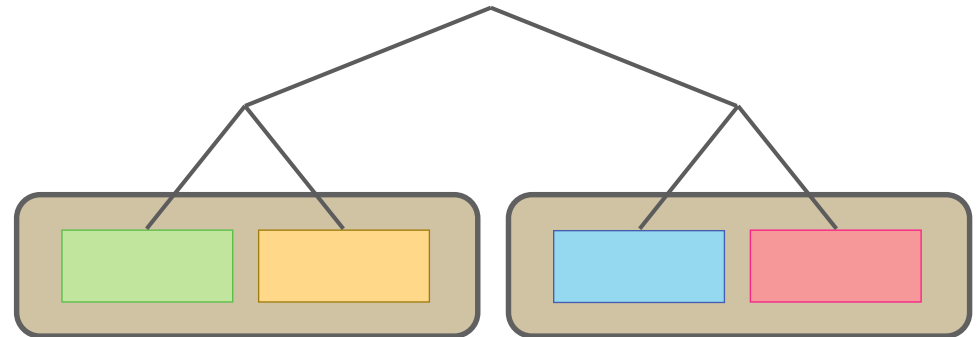
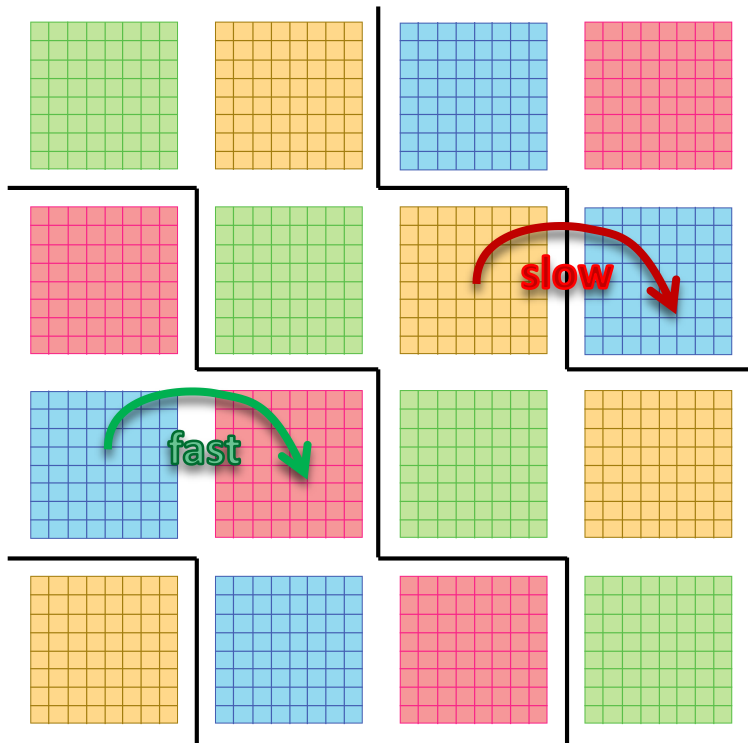


# slow boundaries in halo exchange: 12

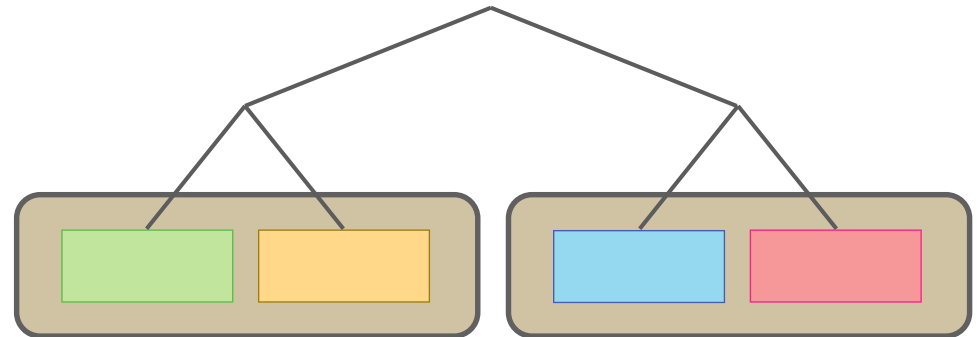
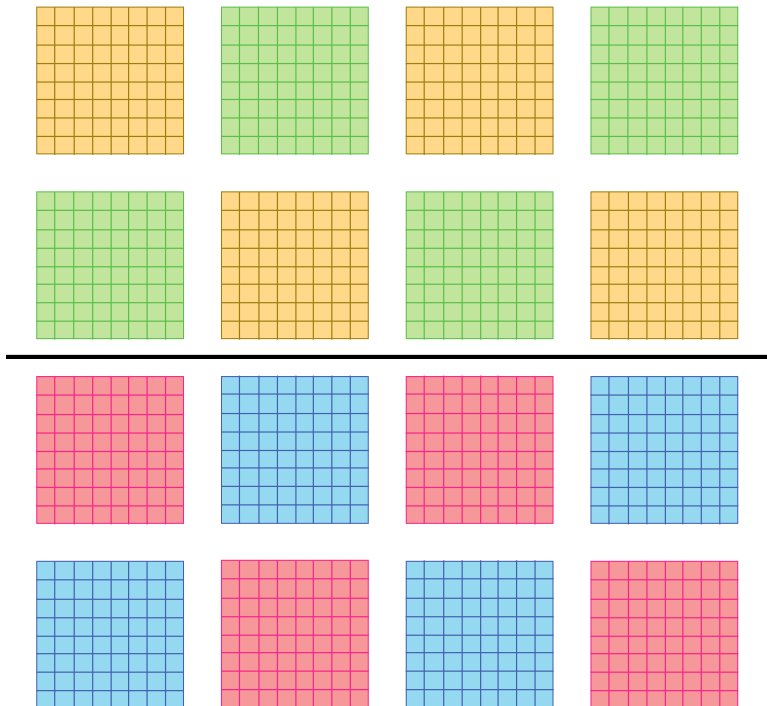


# slow boundaries in halo exchange: 12

# distinct neighbor ranks per process: 2

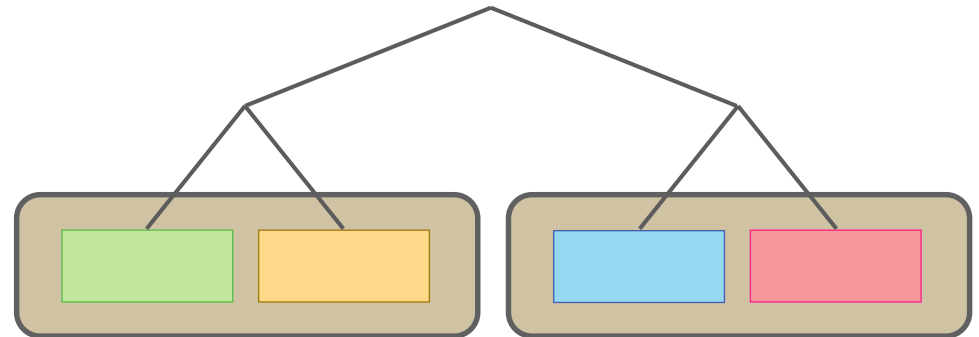
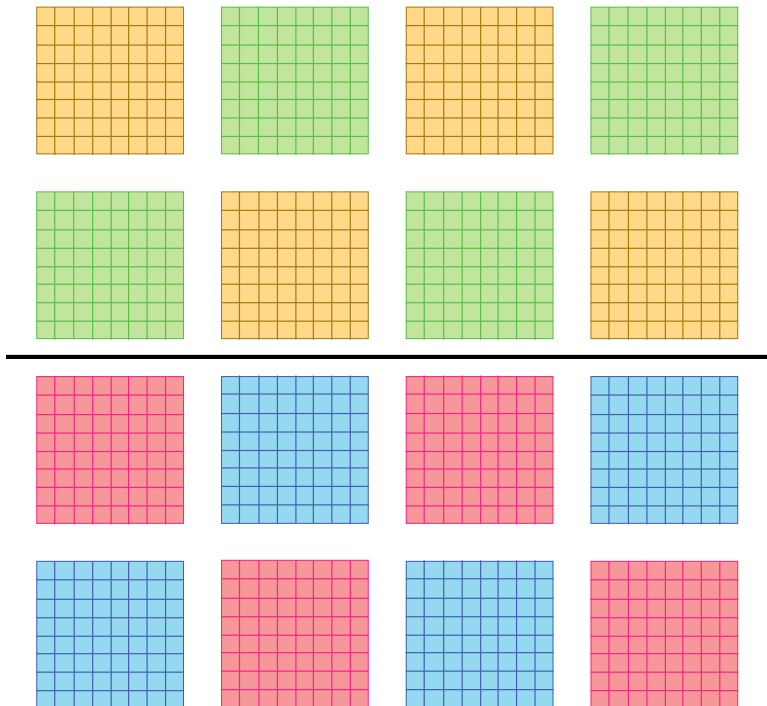


# slow boundaries in halo exchange: 4

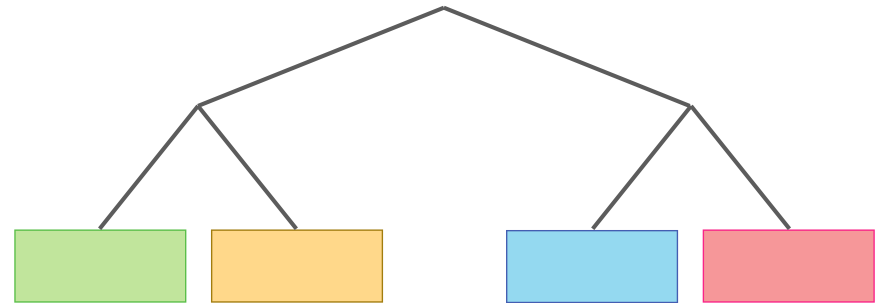
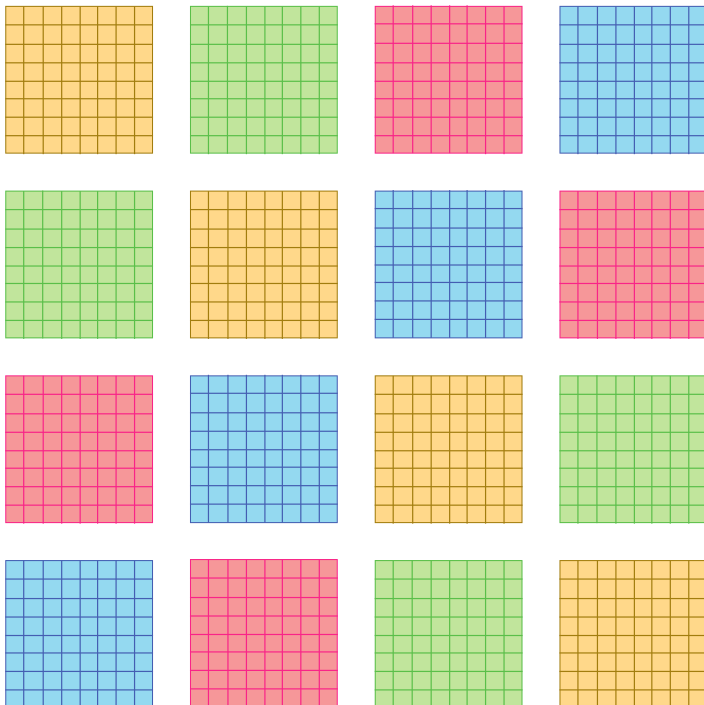


... but isn't time to completion the same?

Time spent in communication per iteration = max. transfer time

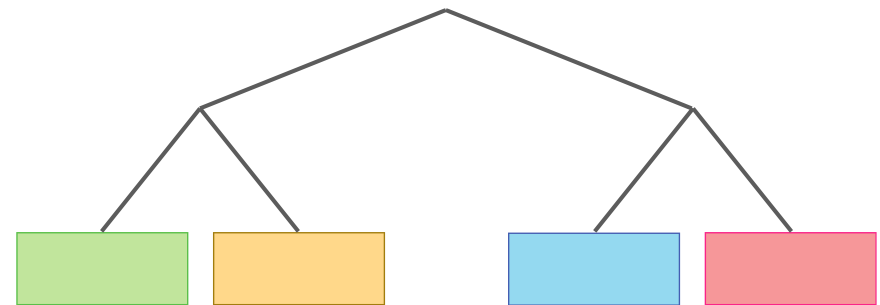
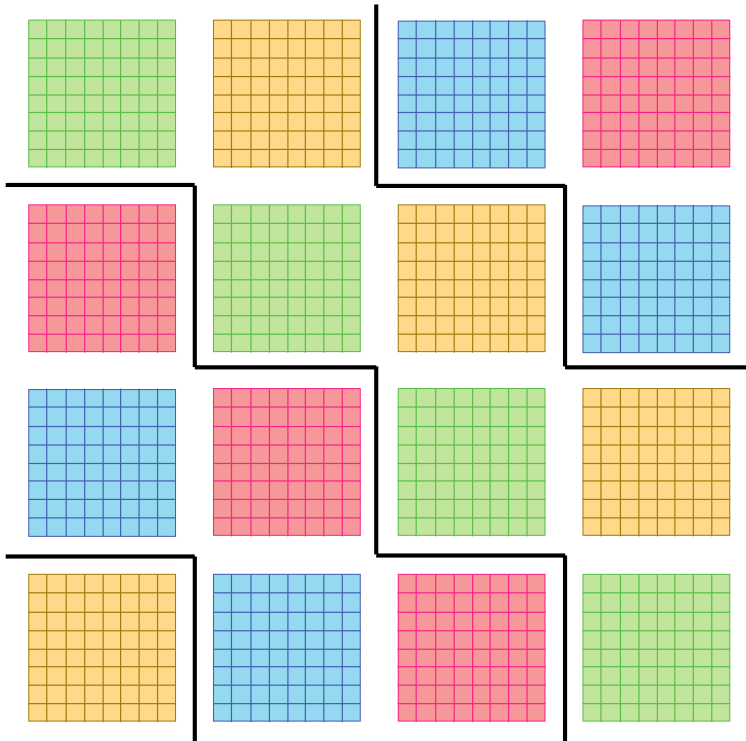


How many slow boundaries for this mapping scheme?

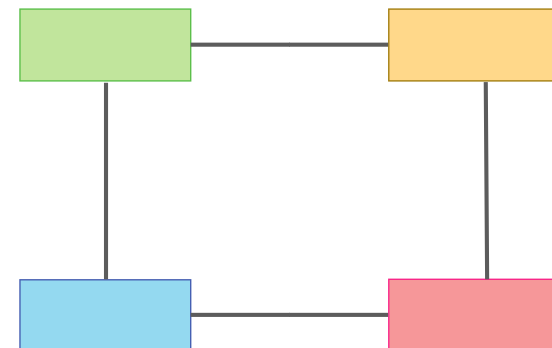




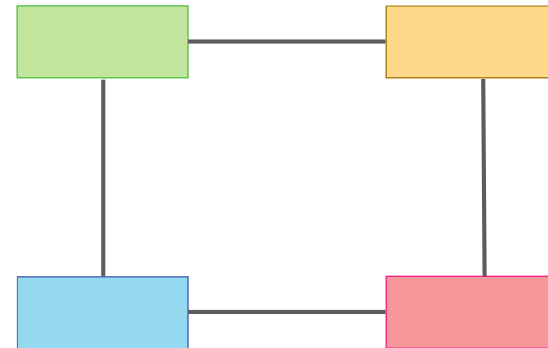
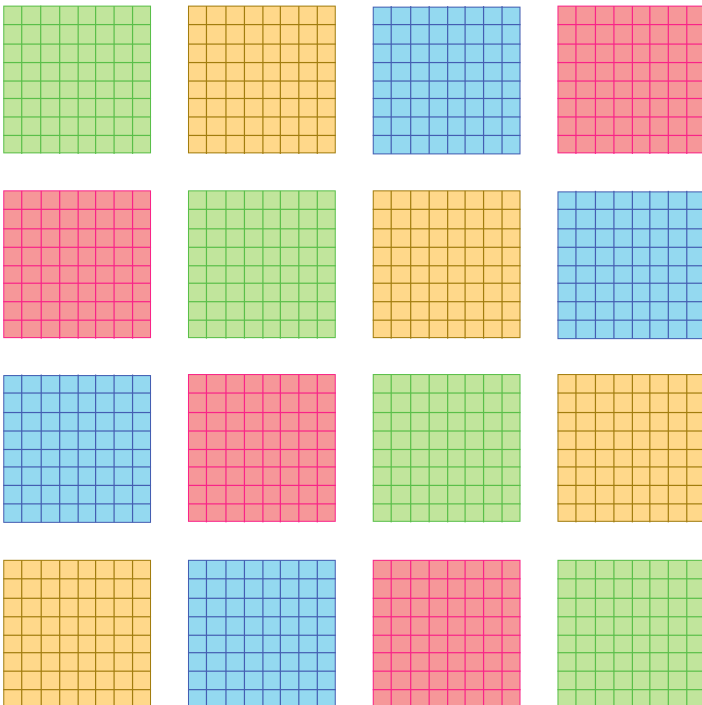
Would a more interconnected, flat hardware topology help?



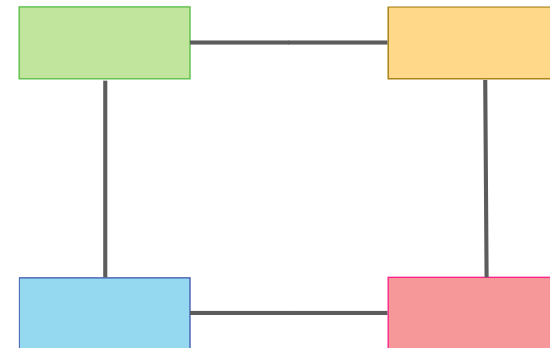
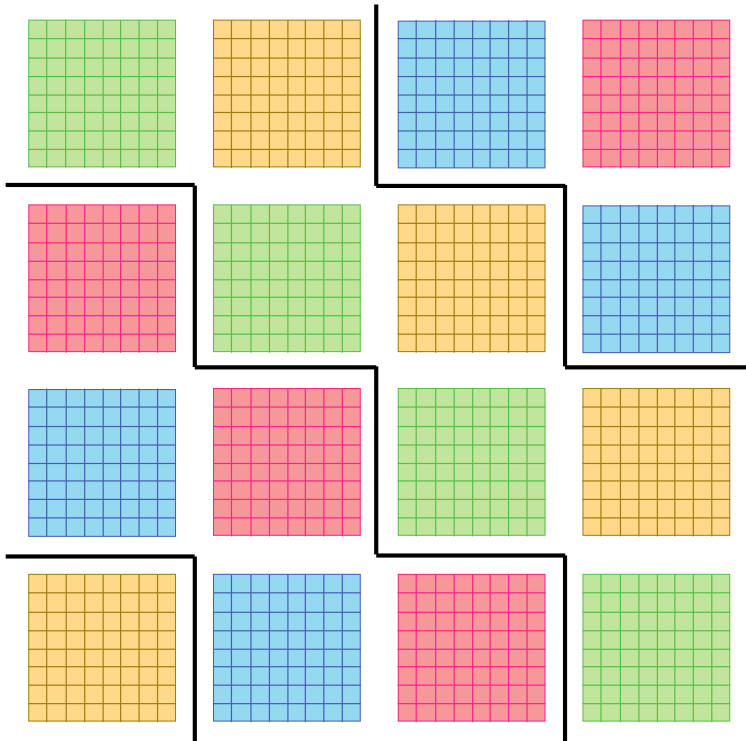
vs.



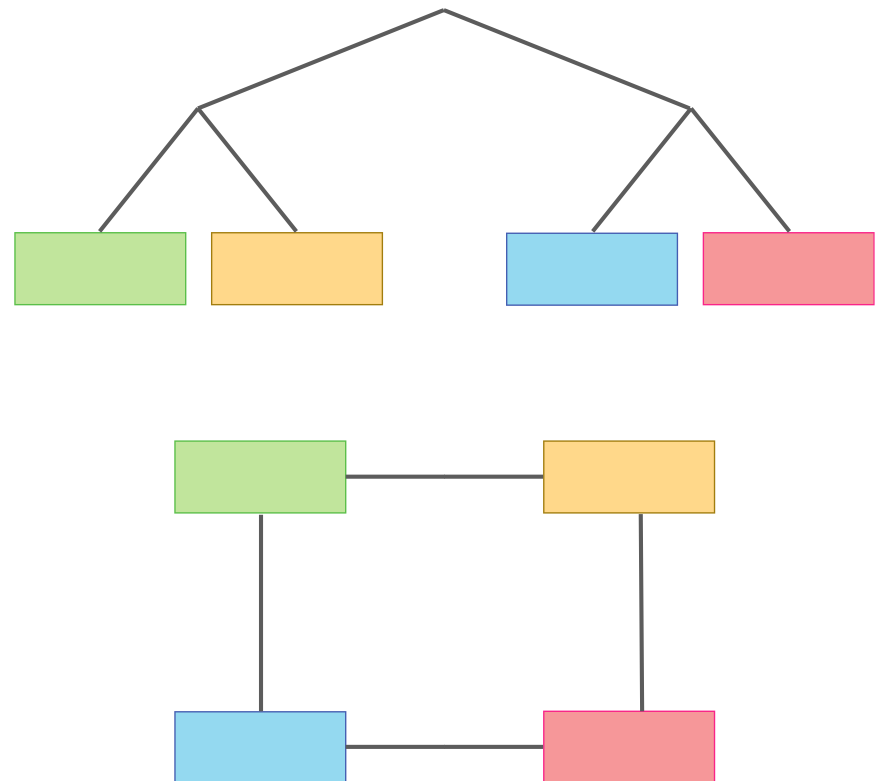
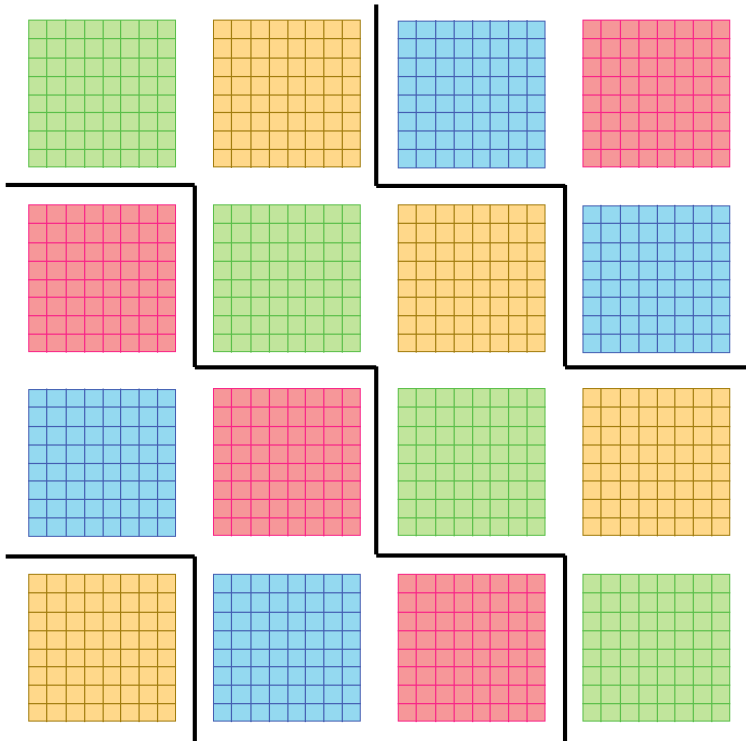
It should!



It should! But domain decomposition scheme prevents to exploit it here.



Just for fun, slow boundaries are even identical for either topology in this example.



## The one MPI tutorial you all want to read:

Basics: <https://cvw.cac.cornell.edu/MPI/>  
P2P: <https://cvw.cac.cornell.edu/MPIP2P/>  
RMA: <https://cvw.cac.cornell.edu/MPloneSided/>  
Advanced: <https://cvw.cac.cornell.edu/MPIAdvTopics/>

Official MPI 3.1 documentation (Index):

<http://www.mpi-forum.org/docs/mpi-3.1/mpi31-report/mpi31-report.htm#Node0>

Again, a collection of documented MPI examples:

<http://www.mcs.anl.gov/~thakur/sc14-mpi-tutorial/>



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