



# Introduction to MPI-I/O

ATPESC 2019

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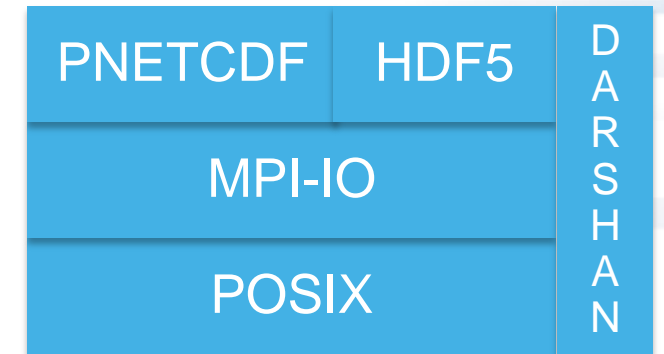
Argonne National Laboratory

Q Center, St. Charles, IL (USA)

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# Plan of attack

- Bottom-up tour of I/O interfaces
  - POSIX routines called by MPI-IO implementations
  - Parallel-NetCDF routines build on top of MPI-IO
- Simple toy programs
  - Refining example several times throughout day
  - We can apply these lessons to your own code in evening session
- Demonstrating some tools for understanding what's going on
- “Game of Life” for your reference



materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Hands on materials

- Code for this ...
  - Simple array I/O
- ... and other sections available on our gitlab site:
  - Game of Life I/O
  - Darshan
  - Burst buffers
  - Globus
- <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>
- I'm going to give you a few minutes to try each hands-on. Can continue working in evening session if you need more time.

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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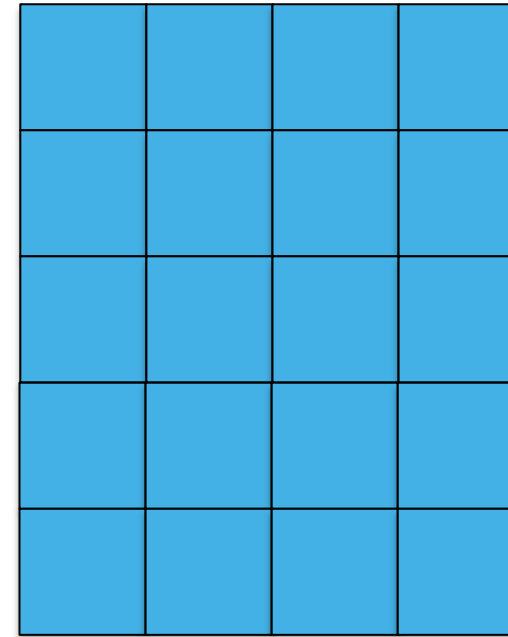
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# Operating on Arrays

- Arrays show up in many scientific applications
  - Matrix operations
  - Particle maps
  - Regions of space
  - Time series
  - Images
- Probably your real application more complicated but an array or two (or more) is in there somewhere, I'd wager.



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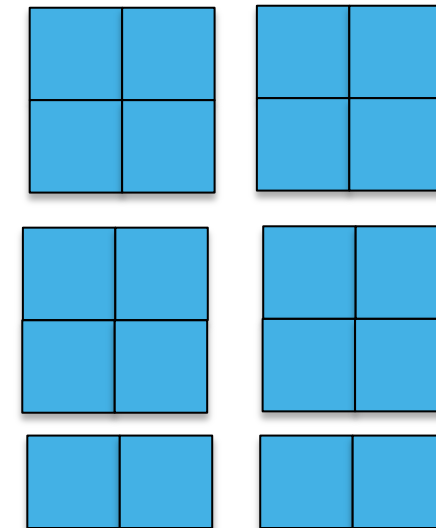
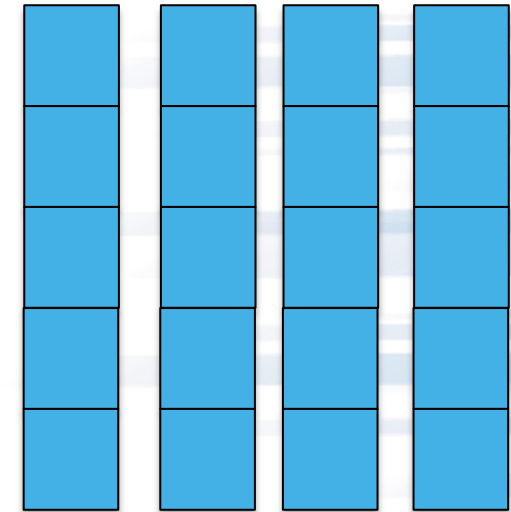
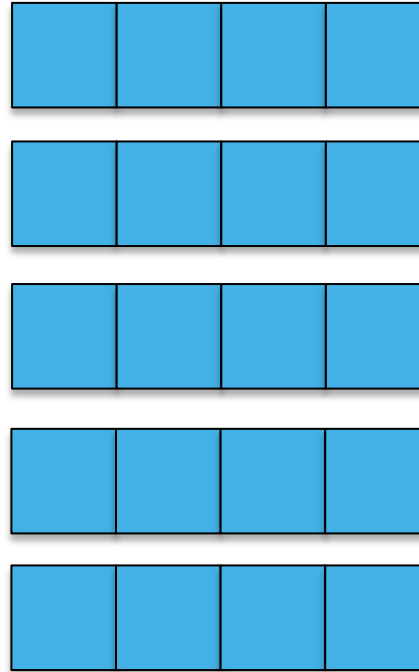
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# Decomposition

- How do we physically access locally parts of a logically larger distributed array in parallel...
  - Piecewise?
  - Chunks?
  - Rows?
- Largely dictated by application algorithm needs
  - E.g. volume rendering math requires chunks not rows.
- Choice impacts memory and I/O performance



materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Supporting Checkpoint/Restart

- For long-running applications, the cautious user checkpoints
- Application-level checkpoint involves the application saving its own state
  - With a bit of extra effort, can be portable
- A canonical representation is preferred
  - Independent of number of processes
- Restarting is then possible
  - Canonical representation aids restarting with a different number of processes
- Also eases data analysis (when using same output)

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Defining a Checkpoint

- Need enough to restart
  - Header information
    - Size of problem (e.g. matrix dimensions)
    - Description of environment (e.g. input parameters)
  - Program state
    - Should represent the global (canonical) view of the data
- Ideally stored in a convenient container
  - Single “thing” (file, object, keyval store...)
- If all processes checkpoint at once, naturally a parallel, collective operation

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# HANDS-ON 1: simple data descriptions

- Consider an application that operates on a 2-d array of integers.
  1. Write code declaring a 2-d array of integers
    - Probably want to allocate on heap, not stack
    - Later steps will be easier if you make it a single allocation
  2. Define a data structure describing the experiment
    - E.g. C struct with row, column, iteration
- Use whatever language you like...
  - ... but Phil and I can only be helpful if you use C
- Source “`setup-env.sh`” to load necessary modules
- Could run this on laptop if you want: shouldn't require any libraries

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# HANDS-ON 1 solutions

C struct holding metadata

```
typedef struct {  
    int row;  
    int col;  
    int iter;  
} science;
```

Do this: index into a single big allocation

```
int *array;  
array = malloc(XDIM*YDIM*sizeof(*array));
```

Don't do this: N allocations will be slower and harder to describe

```
/* not MPI-friendly: describing this memory region will require  
 * a more complicated data type description */  
int **annoying;  
annoying = malloc(YDIM*sizeof(*array));  
for (int i=0; i<YDIM; i++)  
    annoying[i] = malloc(XDIM*sizeof(*array));
```

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# POSIX I/O

- POSIX is the IEEE Portable Operating System Interface for Computing Environments
- “POSIX defines a standard way for an application program to obtain basic services from the operating system”
  - Mechanism almost all serial applications use to perform I/O
- POSIX was created when a single computer owned its own file system

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Deficiencies in serial interfaces

POSIX:

```
fd = open("some_file", O_WRONLY|O_CREAT,  
  S_IRUSR|S_IWUSR);  
ret = write(fd, w_data, nbytes);  
ret = lseek(fd, 0, SEEK_SET);  
ret = read(fd, r_data, nbytes);  
ret = close(fd);
```

FORTRAN:

```
OPEN(10, FILE='some_file', &  
  STATUS="replace", &  
  ACCESS="direct", RECL=16);  
WRITE(10, REC=2) 15324  
CLOSE(10);
```

- 
- Typical (serial) I/O calls seen in applications
  - No notion of other processors
  - Primitive (if any) data description methods
  - Tuning limited to open flags
  - No mechanism for data portability
    - Fortran not even portable between compilers

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# HANDS-ON 2: simple I/O

- We haven't talked about MPI-IO or I/O libraries, but we can still checkpoint.
  - Serial I/O, not parallel
- Implement “write\_data”
  - Will create file and fill in data
  - Prototype:
    - `int write_data(char *filename)`
  - Use system calls (`open()`, `write()`, `close()`), not “stdio” calls (`fopen()`, `fwrite()`, `fclose()`): will map more closely to MPI-IO later
  - How will you know it worked?
  - We are going to repeatedly revise `write_data()` (and later `read_data()`) with each exercise
    - Software engineering: hide details

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# RUNNING

- Submit to the 'R.ATPESC2019' queue
- I've provided a 'submit.sh' shell script
  - `qsub -q R.ATPESC2019 submit.sh <program> [filename]`
    - If you don't give [filename], then 'testfile' used.
- Which Theta file system to use?
  - Tried to make scripts do right thing by default
  - Please don't use the NFS-mounted home directory
  - `submit.sh` should already point you to the right lustre directory
- Make a directory for your data
  - `mkdir -p /lus/theta-fs0/projects/ATPESC2019/data-and-io/$USER`

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# Solution fragments:

```
int write_data(char *filename)
{
    science data = {
        .row = YDIM,
        .col = XDIM,
        .iter = 1
    };

    int *array;
    int fd;
    int ret=0;

    array = buffer_create(0, XDIM, YDIM);

    fd = open(filename, O_CREAT|O_WRONLY,
               S_IRUSR|S_IWUSR);
    ret = write(fd, &data, sizeof(data));
    ret = write(fd, array, XDIM*YDIM*sizeof(int));
    ret = close(fd);

    return ret;
}
materials: https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on
```

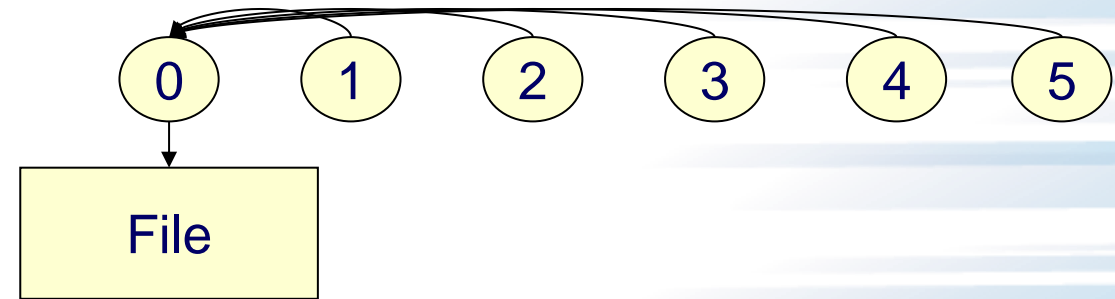
Reading a **binary file**: “cat” won’t work. Could write a c program to read. Several utilities available. I like ‘od’: (historically it only did an “octal dump”). The (t)ype argument can select (d)ecimal

```
% od -td testfile
0000000          1          5          1          0
0000020          1          2          3          4
0000040
```



# HANDS-ON 3: send-to-master

- Parallel program, but serial I/O
  1. `Write_data()` should take an MPI Communicator
  2. Call `MPI_Init()` and `MPI_Finalize()`
  3. Use `MPI_Gather` to collect all data onto rank 0:
- Only rank 0 does I/O; writes header and all array information
- What's good about send-to-master? What's bad?



Hdr			
0	1	2	3
10	11	12	13
20	21	22	23
30	31	32	33
40	41	42	43

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# Solution fragments: MPI\_Gather, write larger data from rank 0

```
MPI_Comm_rank(comm, &rank);
MPI_Comm_size(comm, &nprocs);
/* every process creates its own buffer */
array = buffer_create(rank, XDIM, YDIM);

/* and then sends it to rank 0 */
int *buffer =
    malloc(XDIM*YDIM*nprocs*sizeof(int));

MPI_CHECK(MPI_Gather(
    /* sender (buffer,count,type) tuple */
    array, XDIM*YDIM, MPI_INT,
    /* receiver tuple */
    buffer, XDIM*YDIM, MPI_INT,
    /* who gathers and across which context */
    0, comm));
```

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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# Solution fragments: writing from rank 0

```
if (rank == 0) {  
    /* looks like serial with more data */  
    ...  
    /* writing a global array, not just our  
    local piece of it */  
    data.row = YDIM*nprocs;  
    data.col = XDIM;  
    data.iter = 1;  
  
    ret = write(fd, &data, sizeof(data));  
    ret = write(fd, buffer,  
                XDIM*YDIM*nprocs*sizeof(int));  
  
    ret = close(fd);  
    return ret;  
}
```

## Other questions:

- Lots of machines (your laptop; Theta) represent integers as 32 bit little endian. What if you ran this code on Mira?
- We wrote row-wise. What if you wanted to write a column of data?
- What impact would a header have on data layout? Are there other options?

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# HANDS-ON 4: using Darshan

1. Find the darshan log for the last exercise
  2. View the raw counters with “darshan-parser”
  3. Generate a report
    - You might have to transfer PDF locally to view
  4. Find the darshan log for the exercise #2
    - Hint: you can't! – why not?
- Hint: <https://www.alcf.anl.gov/user-guides/darshan>

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Parallel I/O and MPI

- The stdio checkpoint routine works but is not parallel
  - One process is responsible for all I/O
  - Wouldn't want to use this approach for real
- How can we get the full benefit of a parallel file system?
  - We first look at how parallel I/O works in MPI
  - We then implement a fully parallel checkpoint routine
- MPI is a good setting for parallel I/O
  - Writing is like sending and reading is like receiving
  - Any parallel I/O system will need:
    - collective operations
    - user-defined datatypes to describe both memory and file layout
    - communicators to separate application-level message passing from I/O-related message passing
    - non-blocking operations
  - i.e., lots of MPI-like machinery

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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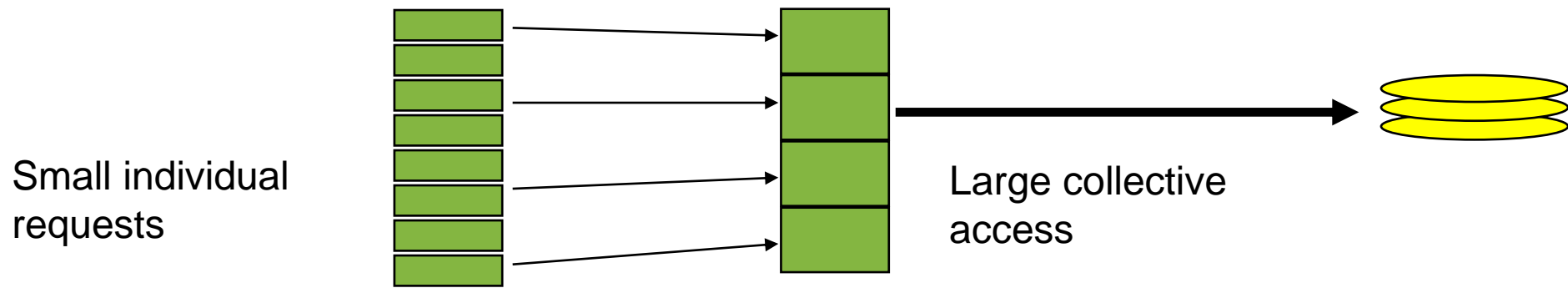
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# Collective I/O

- A critical optimization in parallel I/O
- All processes (in the communicator) must call the collective I/O function
- Allows communication of “big picture” to file system
  - Framework for I/O transformations/optimizations at the MPI-IO layer
  - e.g., two-phase I/O



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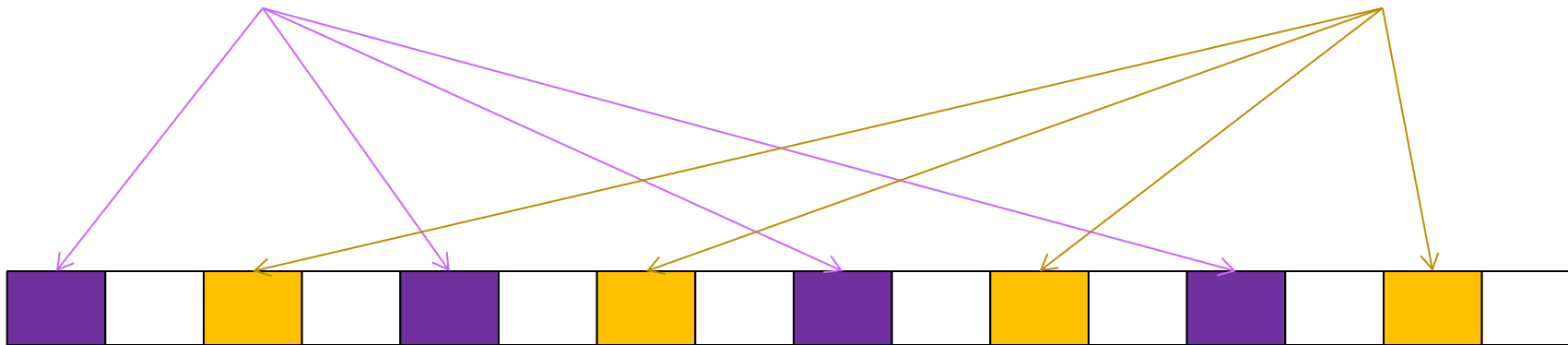
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# Simple MPI-IO

- Collective open: all processes in communicator
- File-side data layout with *file views*
- Memory-side data layout with *MPI datatype* passed to write

```
MPI_File_open(COMM, name, mode,  
             info, fh);  
MPI_File_set_view(fh, disp, etype,  
                filetype, datarep, info);  
MPI_File_write_all(fh, buf, count,  
                 datatype, status);
```

```
MPI_File_open(COMM, name, mode,  
             info, fh);  
MPI_File_set_view(fh, disp, etype,  
                filetype, datarep, info);  
MPI_File_write_all(fh, buf, count,  
                 datatype, status);
```



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# Collective MPI I/O Functions

- Not going to go through the MPI-IO API in excruciating detail
  - Happy to discuss during exercises, evening
- **MPI\_File\_write\_at\_all**, etc.
  - **\_all** indicates that all processes in the group specified by the communicator passed to MPI\_File\_open will call this function
  - **\_at** indicates that the position in the file is specified as part of the call; this provides thread-safety and clearer code than using a separate “seek” call
- Each process specifies only its own access information
  - the argument list is the same as for the non-collective functions
  - OK to participate with zero data
    - All processes must call a collective
    - Process providing zero data might participate behind the scenes anyway

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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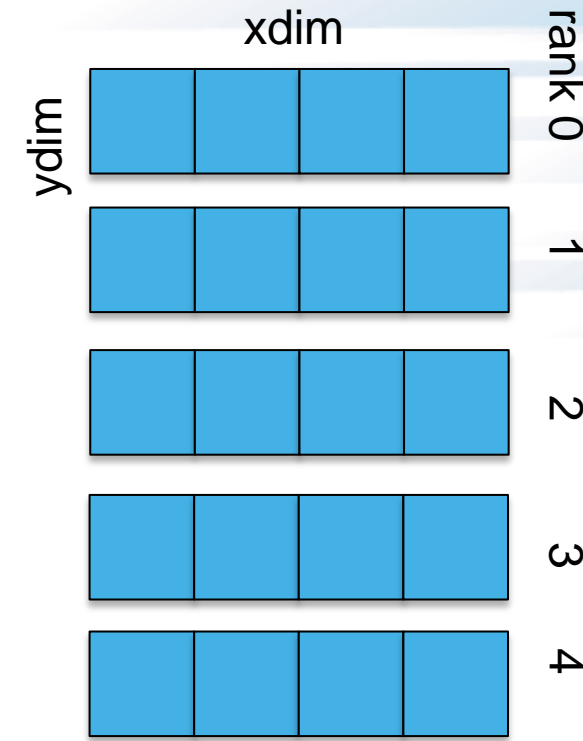
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# HANDS-ON 5: writing with MPI-IO

- Let's take "I/O from master" example and make it parallel
- Use `MPI_File_open` instead of `open`
- Only one process needs to write header
  - Independent `MPI_File_write`
- Every process sets a "file view"
  - Need to skip over header – file view has an "offset" field just for this case
  - The "file view" here is not complicated but we are operating on integers, not bytes:
    - `MPI_File_set_view(fh, sizeof(header), MPI_INT, MPI_INT, "native", info);`
- Each process writes one slice/row of array
  - `MPI_File_write_at_all`
  - Offset "`rank*XDIM*YDIM`"
  - "(buffer, count, datatype)" tuple: `(values, XDIM*YDIM, MPI_INT)`



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# Solution fragments for Hands-On 5

Header I/O from rank 0:

```
if (rank == 0) {  
    MPI_CHECK(MPI_File_write(fh,  
        &header, sizeof(header), MPI_BYTE,  
        MPI_STATUS_IGNORE) );  
}
```

Collective I/O from all ranks

```
MPI_File_write_at_all(fh, rank*XDIM*YDIM,  
    values, XDIM*YDIM, MPI_INT,  
    MPI_STATUS_IGNORE);
```

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# Hands-on 5 continued: Darshan

- A lot like #4: let's use Darshan
  - Find Darshan log file, but don't generate report right away
- What do you think the report will say?
- OK, now generate the report. Were you surprised?

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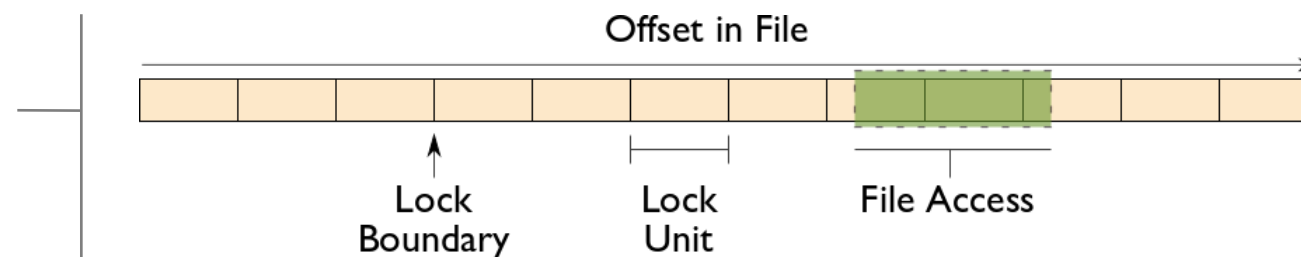


# Managing Concurrent Access

**Files are treated like global shared memory regions. Locks are used to manage concurrent access:**

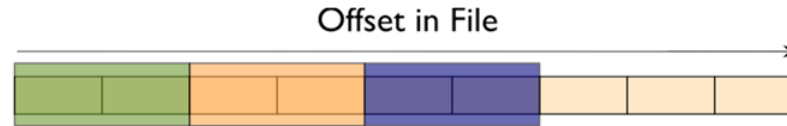
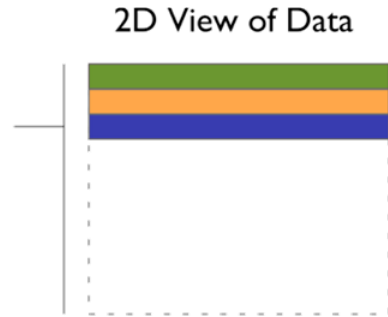
- Files are broken up into lock units
  - Unit boundaries are dictated by the storage system, regardless of access pattern
- Clients obtain locks on units that they will access before I/O occurs
- Enables caching on clients as well (as long as client has a lock, it knows its cached data is valid)
- Locks are reclaimed from clients when others desire access

If an access touches any data in a lock unit, the lock for that region must be obtained before access occurs.



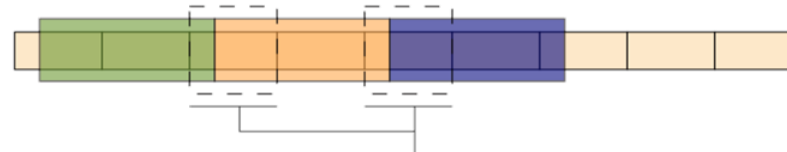
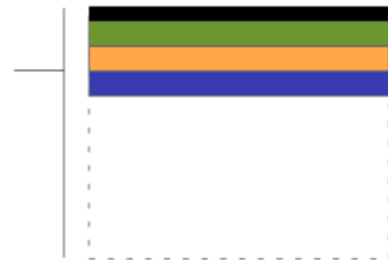
# Implications of Locking in Concurrent Access

The left diagram shows a row-block distribution of data for three processes. On the right we see how these accesses map onto locking units in the file.



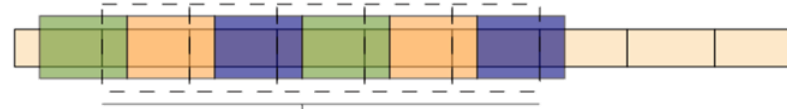
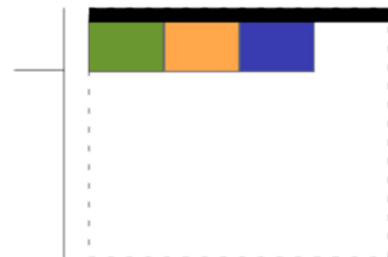
When accesses are to large contiguous regions, and aligned with lock boundaries, locking overhead is minimal.

In this example a header (black) has been prepended to the data. If the header is not aligned with lock boundaries, false sharing will occur.



These two regions exhibit *false sharing*: no bytes are accessed by both processes, but because each block is accessed by more than one process, there is contention for locks.

In this example, processes exhibit a block-block access pattern (e.g. accessing a subarray). This results in many interleaved accesses in the file.

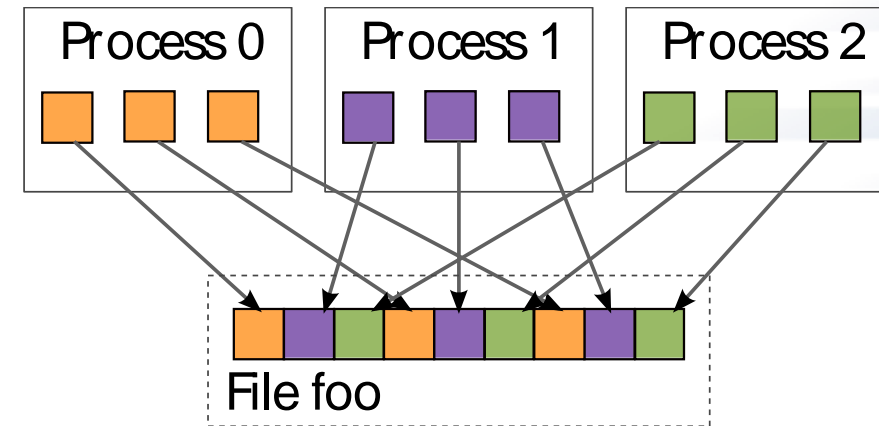


When a block distribution is used, sub-rows cause a higher degree of false sharing, especially if data is not aligned with lock boundaries.

# I/O Transformations

Software between the application and the file system performs transformations, primarily to improve performance.

- Goals of transformations:
  - Reduce number of operations to PFS (avoiding latency)
  - Avoid lock contention (increasing level of concurrency)
  - Hide number of clients (more on this later)
- With “transparent” transformations, data ends up in the same locations in the file as it would have been normally
  - i.e., the file system is still aware of the actual data organization



When we think about I/O transformations, we consider the mapping of data between application processes and locations in file.

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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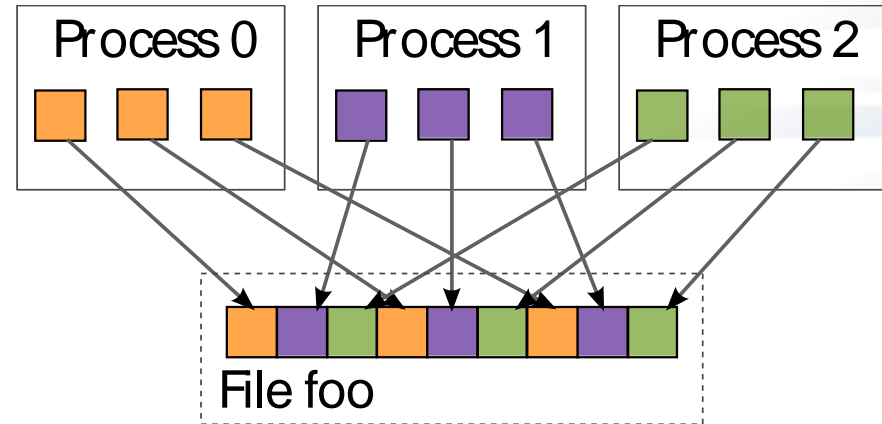


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# I/O Transformations

Software between the application and the file system performs transformations, primarily to improve performance.

- We will tour through a few examples of data transformations in the following slides
- The important thing to remember is that software already exists to do these things for you in HDF5, PnetCDF, ADIOS, and MPI-IO
- If you find yourself replicating these optimizations by hand, look around to see if you can find an off-the-shelf solution

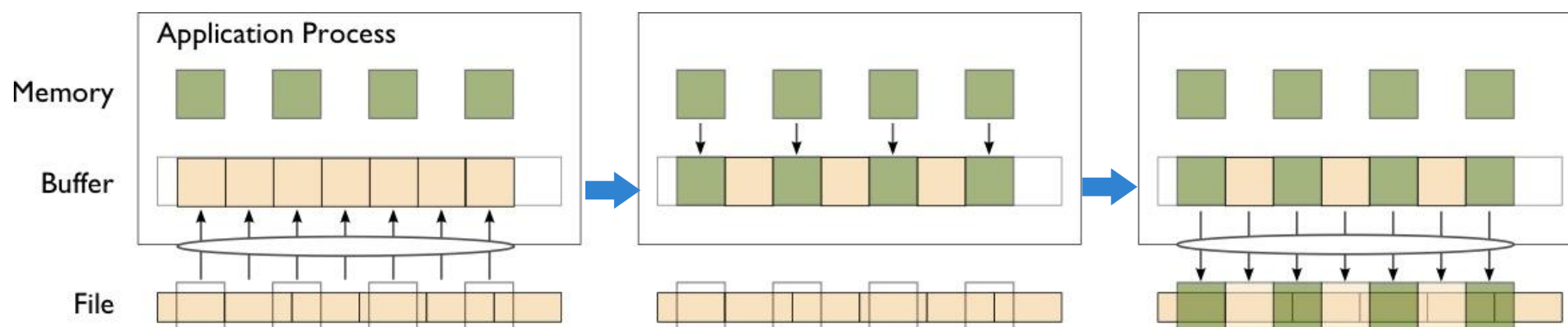


When we think about I/O transformations, we consider the mapping of data between application processes and locations in file.

# Reducing Number of Operations

Because most operations go over multiple networks, I/O to a PFS incurs more latency than with a local FS. *Data sieving* is a technique to address I/O latency by combining operations:

- When reading, application process reads a large region holding all needed data and pulls out what is needed
- When writing, three steps required (below)
- Somewhat counter-intuitive: do extra I/O to avoid contention



**Step 1:** Data in region to be modified are read into intermediate buffer (1 read).

**Step 2:** Elements to be written to file are replaced in intermediate buffer.

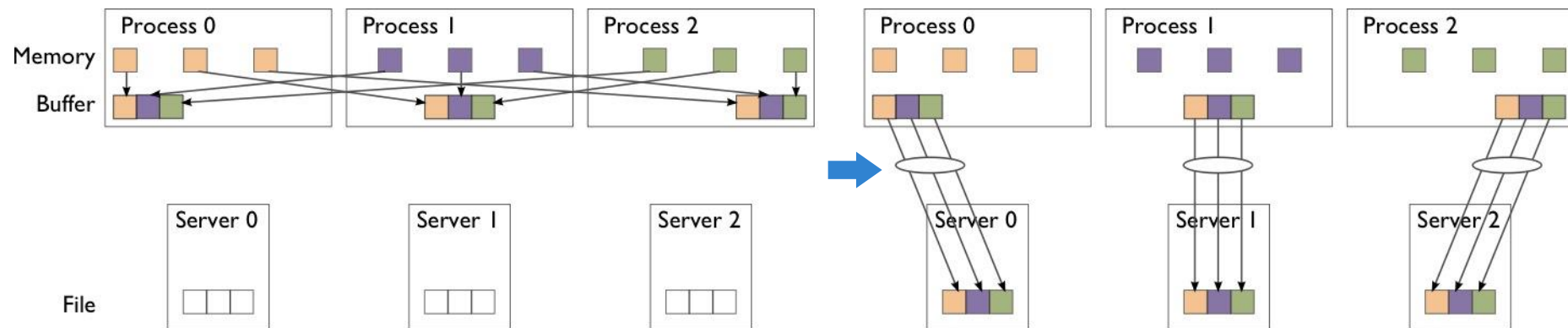
**Step 3:** Entire region is written back to storage with a single write operation.

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>

# Avoiding Lock Contention

**We can reorder data among processes to avoid lock contention.** *Two-phase I/O* splits I/O into a data reorganization phase and an interaction with the storage system (two-phase write depicted):

- Data exchanged between processes to match file layout
- 0<sup>th</sup> phase determines exchange schedule (not shown)



**Phase 1:** Data are exchanged between processes based on organization of data in file.

**Phase 2:** Data are written to file (storage servers) with large writes, no contention.

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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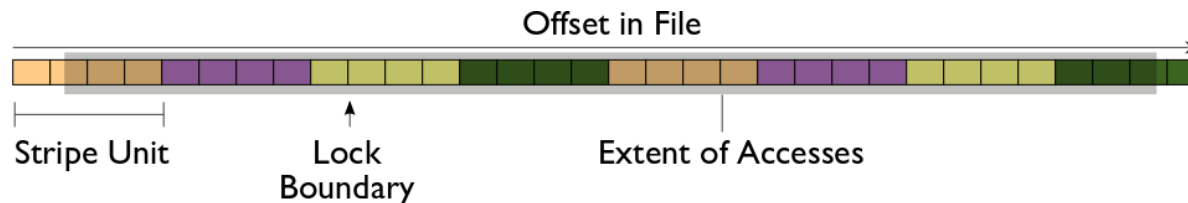


# Two-Phase I/O Algorithms

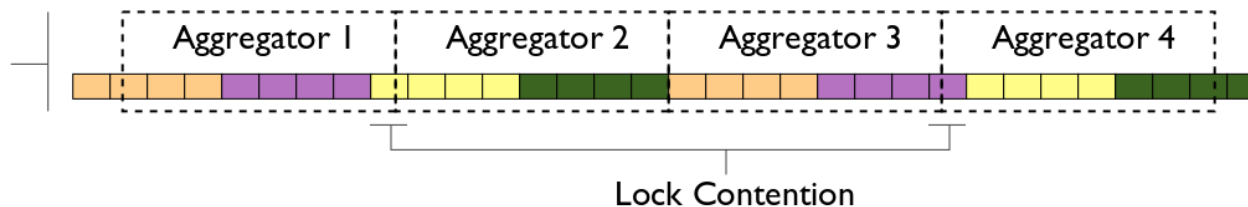
(or, You don't want to do this yourself...)

For more information, see W.K. Liao and A. Choudhary, "Dynamically Adapting File Domain Partitioning Methods for Collective I/O Based on Underlying Parallel File System Locking Protocols," SC2008, November, 2008.

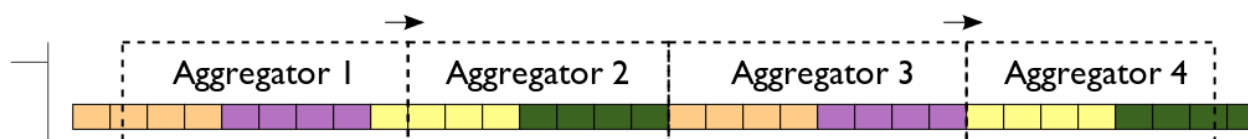
Imagine a collective I/O access using four aggregators to a file striped over four file servers (indicated by colors):



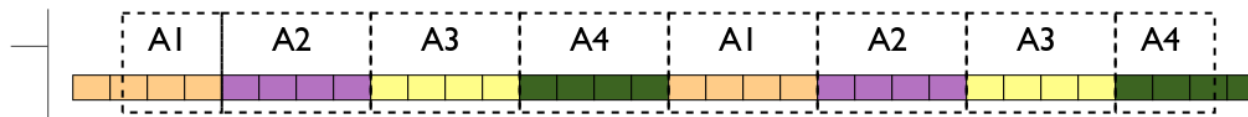
One approach is to evenly divide the region accessed across aggregators.



Aligning regions with lock boundaries eliminates lock contention.



Mapping aggregators to servers reduces the number of concurrent operations on a single server and can be helpful when locks are handed out on a per-server basis (e.g., Lustre).

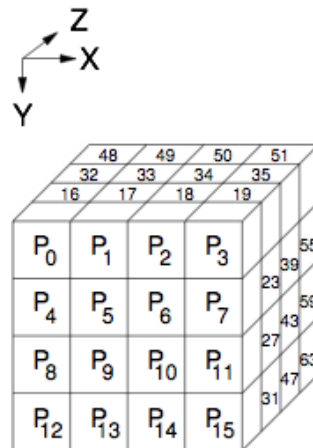
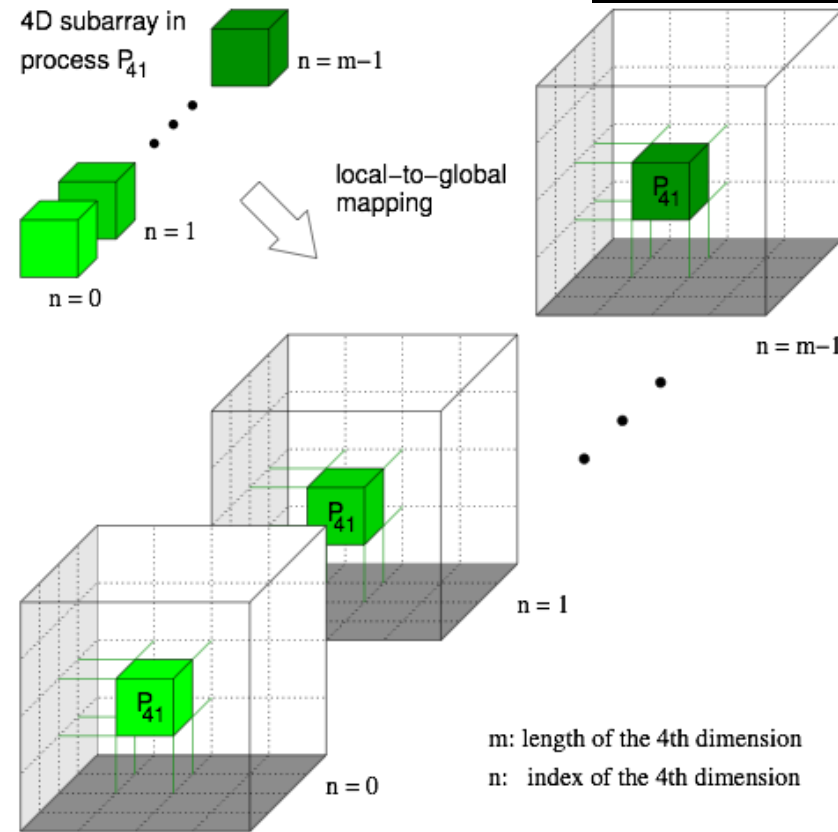
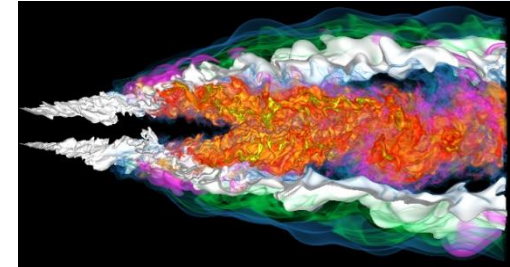


Today's systems also choose aggregators that are "best" for storage

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>

# S3D Turbulent Combustion Code

- S3D is a turbulent combustion application using a direct numerical simulation solver from Sandia National Laboratory
- Checkpoints consist of four global arrays
  - 2 3-dimensional
  - 2 4-dimensional
  - 50x50x50 fixed subarrays



Thanks to Jackie Chen (SNL), Ray Grout (SNL), and Wei-Keng Liao (NWU) for providing the S3D I/O benchmark, Wei-Keng Liao for providing this diagram, C. Wang, H. Yu, and K.-L. Ma of UC Davis for image.

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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# Impact of Transformations on S3D I/O

- Testing with PnetCDF output to single file, three configurations, 16 processes
  - All MPI-IO optimizations (collective buffering and data sieving) disabled
  - Independent I/O optimization (data sieving) enabled
  - Collective I/O optimization (collective buffering, a.k.a. two-phase I/O) enabled

Application did the same thing in every case

	Coll. Buffering and Data Sieving Disabled	Data Sieving Enabled	Coll. Buffering Enabled (including Aggregation)
POSIX writes	102,401	81	<b>5</b>
POSIX reads	0	80	0
MPI-IO writes	64	64	64
Unaligned in file	102,399	80	4
Total written (MB)	6.25	<b>87.11</b>	6.25
Runtime (sec)	1443	11	6.0
Avg. MPI-IO time per proc (sec)	<b>1426.47</b>	4.82	0.60

materials: <https://xgtilab.cels.anl.gov/ATPESC-IO/hands-on>



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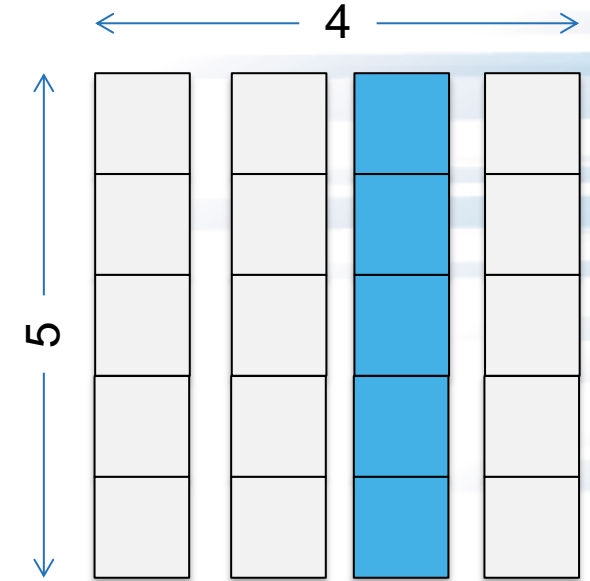
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# HANDS-ON 6: reading with MPI-IO

- Slightly different: all processes read one row
  - For simplicity, same row
- File view will be more complicated, use MPI “Subarray” datatype
- In C, array access is described in “row-major”
  - `array_size[0] = 5; array_size[1] = 4;`
- File view uses derived ‘subarray’, not built-in MPI\_INT
- Location in file given with subarray type; no offset in `MPI_File_read_all`
  - Still provide a “buffer, count, datatype” tuple for memory layout



# Solution fragments

## Type creation

```
/* In C-order the arrays are row-major:
 *
 * |-----|
 * |-----|
 * |-----|
 *
 * The 'sizes' of the above array would be 3,5
 * The last column would be a "subsize" of 3,1
 * And a "start" of 0,5 */

sizes[0] = nprocs; sizes[1] = XDIM;
sub[0] = nprocs; sub[1] = 1;
starts[0] = 0; starts[1] = XDIM/2;

MPI_Type_create_subarray(NDIMS,
    sizes, sub, starts,
    MPI_ORDER_C, MPI_INT, &subarray);
MPI_Type_commit(&subarray);
```

## File view and read

```
MPI_CHECK(MPI_File_set_view(fh, sizeof(header),
    MPI_INT, subarray, "native", info));
MPI_Type_free(&subarray);
MPI_CHECK(MPI_File_read_all(fh,
    read_buf, nprocs, MPI_INT, MPI_STATUS_IGNORE);
```

# Hands on 6 continued: Darshan

- How does this workload differ from the write?
- Change the 'read\_all' to an independent 'read'
  - What do you think the Darshan output will say? Find out.

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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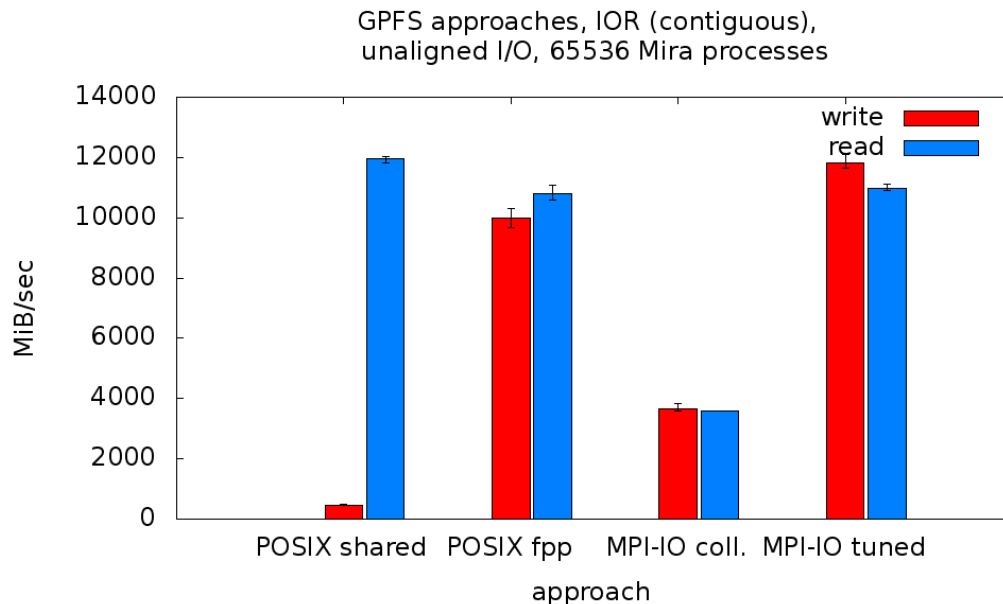
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# GPFS Access three ways

- POSIX shared vs MPI-IO collective
  - Locking overhead for unaligned writes hits POSIX hard
- Default MPI-IO parameters not ideal
  - Reported to IBM; simple tuning brings MPI-IO back to parity
  - “Vendor Defaults” might give you bad first impression
- File per process (fpp) extremely seductive, but entirely untenable on current generation.



materials: <https://xgitylab.cels.anl.gov/ATPESC-IO/hands-on>



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# MPI-IO Takeaway

- Sometimes it makes sense to build a custom library that uses MPI-IO (or maybe even MPI + POSIX) to write a custom format
  - e.g., a data format for your domain already exists, need parallel API
- We've only touched on the API here
  - There is support for data that is noncontiguous in file and memory
  - There are independent calls that allow processes to operate without coordination
- In general we suggest using data model libraries
  - They do more for you
  - Performance can be competitive

materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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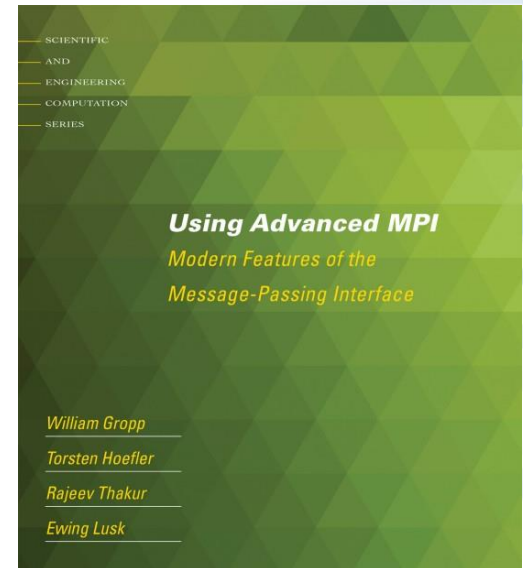


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# MPI-IO References

- On Cray systems, “man intro\_mpi” for 3,000 lines of tuning parameters, debug configuration
- *Using Advanced MPI*, Gropp, Hoeffler, Thakur, Lusk
  - Chapter on MPI I/O routines covers entire API as well as consistency semantics



materials: <https://xgitlab.cels.anl.gov/ATPESC-IO/hands-on>



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