

Scalable and Modular Parallel I/O for Open MPI

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Outline

- Motivation
- MPI I/O: basic concepts
- OMPIO module and parallel I/O frameworks in Open MPI
- Parallel I/O research
- Conclusions and future work



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Motivation

- Study by LLNL (2005):
 - 1 GB/s I/O bandwidth required per Teraflop compute capability
 - Write to the filesystem dominates reading from it by a factor of 5
- Current High-End Systems:
 - K Computer: ~11 PFLOPS, ~96 GB/s I/O bandwidth using 864 OSTs
 - Jaguar (2010): ~1 PFLOPS, ~90 GB/s I/O bandwidth using 672 OSTs

⇒ Gap between available I/O performance and required I/O performance.

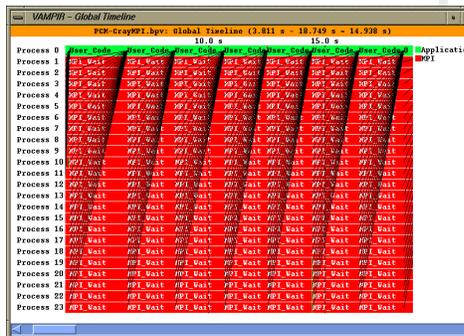


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Application Perspective

- Sequential I/O
 - A single process executes file operations
 - Leads to load imbalance
- Individual I/O
 - Each process has its own files
 - Pre/Post-processing required
- Parallel I/O
 - Multiple processes access (different parts of) the same file (efficiently)



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Part I: MPI I/O



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MPI I/O

- MPI (Message Passing Interface) version 2 introduced the notion of parallel I/O
 - **Collective I/O** : group I/O operations
 - **File view**: registering an access plan to the file in advance
 - **Hints**: application hints on the lanned usage of the
 - **Relaxed consistency semantics**: updates to a file might initially only be visible to the process performing the action
 - **Non-blocking I/O**: asynchronous I/O operations



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General file manipulation functions

```
MPI_File_open ( MPI_Comm comm, char *filename,
                int amode, MPI_Info info,
                MPI_File *fh );
```

- **Collective operation**
 - All processes have to provide the same amode
 - comm must be an intra-communicator
- **Values for amode**
 - MPI_MODE_RDONLY, MPI_MODE_WRONLY, MPI_MODE_RDWR,
 - MPI_MODE_CREATE, MPI_MODE_APPEND, ...
- **Combination of several amodes possible, e.g**
 - **C:** (MPI_MODE_CREATE | MPI_MODE_WRONLY)
 - **Fortran:** MPI_MODE_CREATE + MPI_MODE_WRONLY

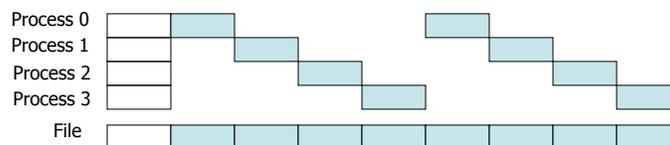


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File View

- **File view: portion of a file visible to a process**
 - Processes can share a common view
 - Views can overlap or be disjoint
 - Views can be changed during runtime
 - A process can have multiple instances of a file open using different file views

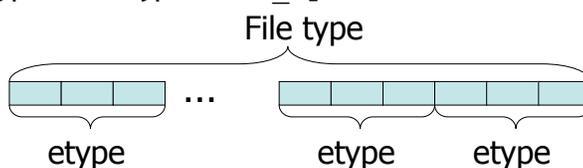


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File View

- Elementary type (etype) : basic unit of the data accessed by the program
- File type: datatype used to construct the file view
 - consists logically of a series of etypes
 - must not have overlapping regions if used in write operations
 - displacements must increase monotonically
- Default file view:
 - displacement = 0
 - etype = file type = MPI_Byte



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Setting a file view

```
MPI_File set_view ( MPI_File fh, MPI_Offset disp,
                   MPI_Datatype etype, MPI_Datatype filetype,
                   char *datarep, MPI_Info info );
```

- The argument list
 - disp: start of the file view
 - etype and filetype: as discussed previously
 - datarep: data representation used
 - info: hints to the MPI library (discussed later)
- Collective operation
 - datarep and extent of etype have to be identical on all processes
 - filetype, disp and info might vary
- Resets file pointers to zero



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File Interoperability

- Fifth parameter of `MPI_File_set_view` sets the data representation used:
 - `native`: data is stored in a file exactly as it is in memory
 - `internal`: data representation for heterogeneous environments using the same MPI I/O implementation
 - `external32`: portable data representation across multiple platforms and MPI I/O libraries.
- User can register its own data representation, providing the according conversion functions (`MPI_Register_datarep`)



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General file manipulation functions

```
MPI_File_read ( MPI_File fh, void *buf, int cnt,
                MPI_Datatype dat, MPI_Status *stat);
MPI_File_write ( MPI_File fh, void *buf, int cnt,
                 MPI_Datatype dat, MPI_Status *stat);
```

- Buffers described by the tuple of
 - Buffer pointer
 - Number of elements
 - Datatype
- Interfaces support data conversion if necessary



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MPI I/O non-collective functions

Positioning	Synchronism	Function
Individual file pointers	Blocking	MPI_File_read
		MPI_File_write
	Non-blocking	MPI_File_iread
		MPI_File_iwrite
Explicit offset	Blocking	MPI_File_read_at
		MPI_File_write_at
	Non-blocking	MPI_File_iread_at
		MPI_File_iwrite_at
Shared file pointers	Blocking	MPI_File_read_shared
		MPI_File_write_shared
	Non-blocking	MPI_File_iread_shared
		MPI_File_iwrite_shared



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Individual I/O in parallel applications

Process 0:

1	2	3	4	5	6	7	8
9	10	11	12	13	14	15	16
17	18	19	20	21	22	23	24
25	26	27	28	29	30	31	32

```
read(..., offset=0, length=2)
read(..., offset=8, length=2)
read(..., offset=16, length=2)
read(..., offset=24, length=2)
```

- Individual Read/Write operations on a joint file often lead to many, small I/O requests from each process
- Arbitrary order of I/O requests from the file system perspective
 - Will lead to suboptimal performance



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Collective I/O in parallel applications

Process 0:

1	2	3	4	5	6	7	8
9	10	11	12	13	14	15	16
17	18	19	20	21	22	23	24
25	26	27	28	29	30	31	32

```
read(..., offset=0, length=4)
MPI_Send (... ,length=2,dest=3,...)
read(..., offset=8, length=4)
MPI_Send (... ,length=2,dest=3,...)
read(..., offset=16, length=4)
MPI_Send (... ,length=2,dest=3,...)
read(..., offset=24, length=4)
MPI_Send (... ,length=2,dest=3,...)
```

- Collective I/O:
 - Offers the potential to rearrange I/O requests across processes, e.g. minimize file pointer movements, minimize locking occurring on the file system level
 - Offers performance benefits if costs of additional data movements < benefit of fewer repositioning of file pointers



Collective I/O: Two-phase I/O algorithm

- Re-organize data across processes to match data layout in file
- Combination of I/O and (MPI level) communication used to read/write data from/to file
- Only a subset of processes actually touch the file (aggregators)
- Large read/write operations split into multiple cycles internally
 - Limits the size of temporary buffers
 - Overlaps communication and I/O operations



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Shared File Pointer Operations

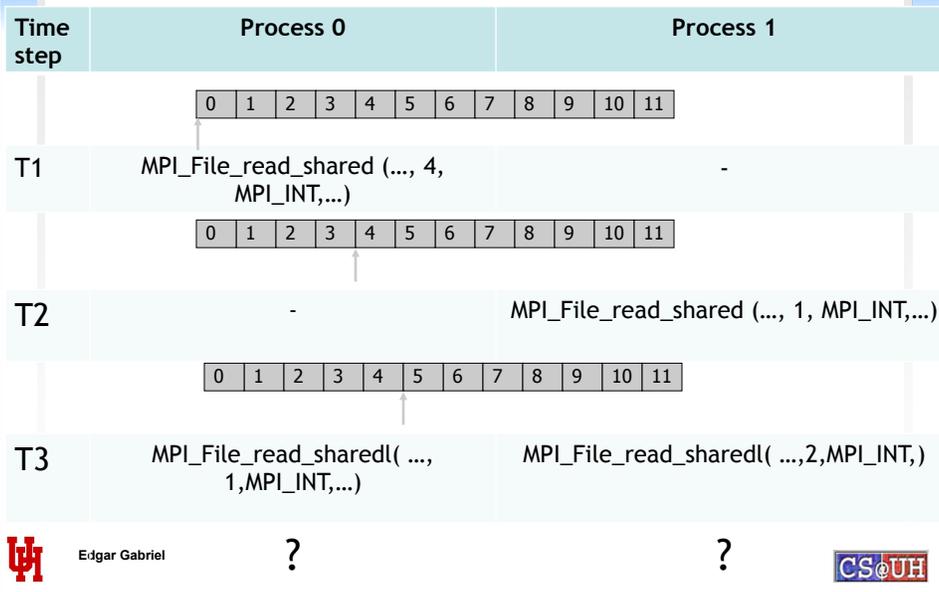
- Shared file pointers: a file pointer shared by a the group of processes that has been used to open the file
 - All processes must have identical file view
 - Might lead to non-deterministic behavior
- Shared file pointer must not interfere with the individual file pointer of each process
- Typical usage scenarios
 - Writing a parallel log-file
 - Work distribution across processes by reading data from a joint file



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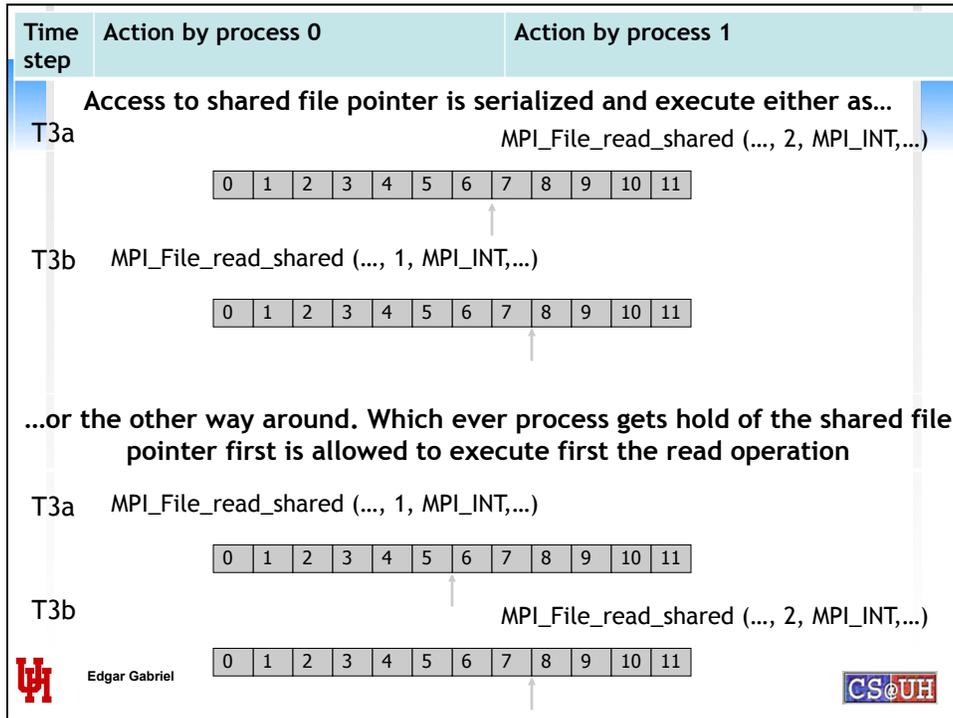


Shared file pointer example



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Consistency of file operation

- MPI does not provide sequential consistency across all processes per default
 - Write on one process is initially just visible on the same process
- Two possibilities to change this behavior

```
MPI_File_set_atomicity ( MPI_File fh, int flag );
```

- If flag = true, all write operations are atomic
- Collective operation

```
MPI_File_sync ( MPI_File fh );
```

- Flushes all write operations on the calling process' file instance

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Hints supported by MPI I/O (I)

Hint	Explanation	Possible values
<code>access_style</code>	Specifies manner in which the file is accessed	<code>read_once</code> , <code>write_once</code> , <code>read_mostly</code> , <code>write_mostly</code> , <code>sequential</code> , <code>reverse_sequential</code> , <code>random</code>
<code>collective_buffering</code>	Use collective buffering ?	<code>true</code> , <code>false</code>
<code>cb_block_size</code>	Block size used for collective buffering	Integer
<code>cb_buffer_size</code>	Total buffer space that can be used for collective buffering	Integer, multiple of <code>cb_block_size</code>
<code>cb_nodes</code>	Number of target nodes used for collective buffering	Integer



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Hints supported by MPI I/O (II)

Hint	Explanation	Possible values
<code>io_node_list</code>	List of I/O nodes that should be used	Comma separated list of strings
<code>nb_proc</code>	Specifies the number of processes typically accessing the file	Integer
<code>num_io_nodes</code>	Number of I/O nodes available in the system	Integer
<code>striping_factor</code>	Number of I/O nodes that should be used for file striping	Integer
<code>striping_unit</code>	Stripe depth	integer



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Part II: OMPIO



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OMPIO Design Goals (I)

- Highly modular architecture for parallel I/O
 - Maximize code reuse, minimize code replication
- Generalize the selection of modules
 - Collective I/O algorithms
 - Shared file pointer operations
- Tighter Integration with Open MPI library
 - Derived data type optimizations
 - Progress engine for non-blocking I/O operations
 - External data representations etc.



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OMPIO Design Goals (II)

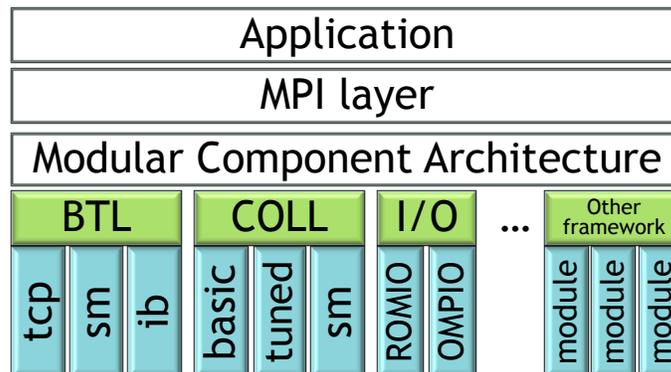
- Adaptability
 - Enormous diversity of I/O hardware and software solutions
 - Number of storage server, bandwidth of each storage server
 - Network connectivity in-between I/O nodes, between compute and I/O nodes, and message passing network between compute nodes
 - Ease the modification of module parameters
 - Ease the development and dropping in of new modules



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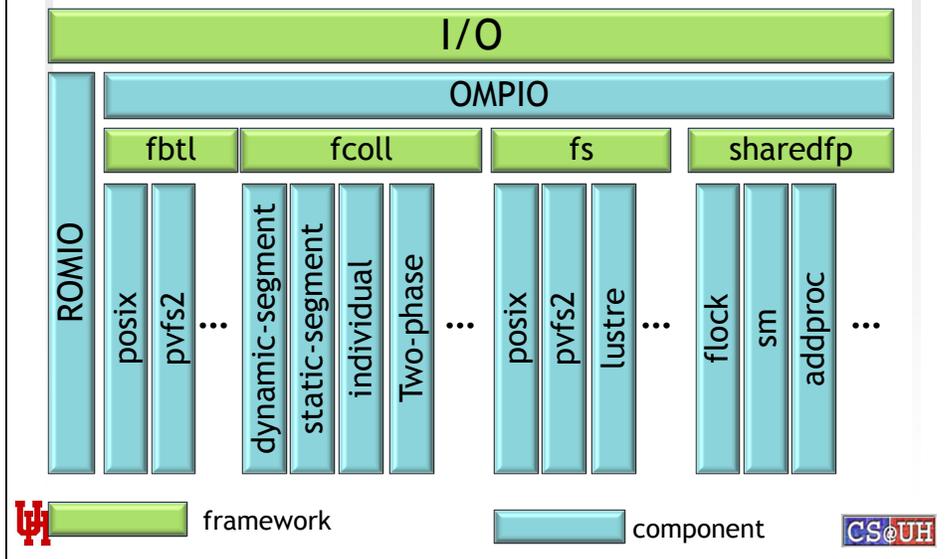
Open MPI Architecture



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OMPIO frameworks overview



OMPIO

- Main I/O component
- 'Understands' MPI semantics
- Translates MPI write/read operations into lower layer operations
- Provides the implementation and the operation of the
 - `MPI_File` handle
 - File view operations
 - (`MPI_Request` structures)
- Triggers upon selection the `fcoll`, `fs`, `fbt1` and `sharedfp` selection logic

fbtl: file byte transfer layer

- Abstraction for individual read and write operations
- A process will have per MPI file one or more fbtl modules loaded
- Main interfaces work with the tuple of `<buffer pointer, length, position in file>`
- Interface:
 - `pwritev()` - `ipwritev()`
 - `preadv()` - `ipreadv()`
 - `progress()`



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fcoll: collective I/O framework

- Provides implementations of the collective I/O operations of the MPI specification
 - `read_all()` - `read_all_begin()/end()`
 - `write_all()` - `write_all_begin()/end()`
 - `read_at_all()` - `read_at_all_begin()/end()`
 - `write_at_all()` - `write_at_all_begin()/end()`
- Selection logic triggered upon setting the file view



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fcoll: selection logic

- Decision between different collective modules based on:
 - `ss`: stripe size of the file system
 - `c`: average contiguous chunk size in file view
 - `k`: minimum data size to saturate write/read bandwidth from one process
 - size of gap in the file view between processes.

Characteristic	Gap Size	Algorithm
$c > k$ and $c > ss$	any	individual
$c \leq k$ and $c > ss$	0	dynamic segmentation
$c < k$ and $c \leq ss$	0	two-phase
$c < k$	> 0	static segmentation



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fs: file system framework

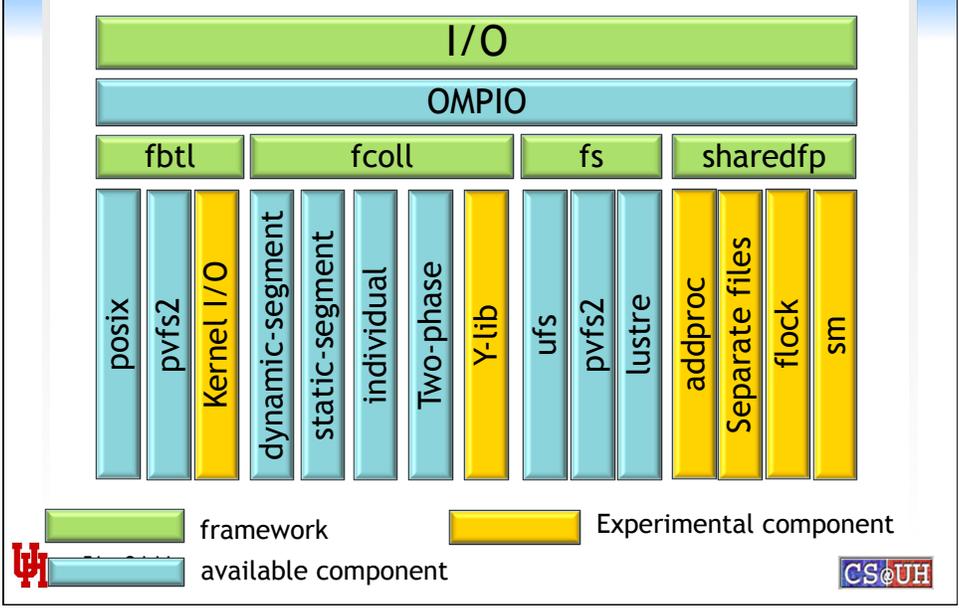
- Handles all file-system related operation
 - Interfaces have mostly collective notion
- Interface:
 - `open()`
 - `close()`
 - `delete()`
 - `sync()`
- Current Lustre and PVFS2 fs components allow to modify stripe size, stripe depth and I/O servers used



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Current status



Performance results: Tile I/O

Shark cluster at University of Houston (PVFS2):

No. of procs.	Tile Size	fcoll module	OMPIO bandwidth	ROMIO bandwidth
81	64 Bytes	Two-phase	591 MB/s	303 MB/s
81	1 MB	Dynamic Segm.	625 MB/s	290 MB/s

Deimos cluster at TU Dresden (Lustre):

No. of procs.	Tile Size	fcoll module	OMPIO bandwidth	ROMIO bandwidth
256	64 Bytes	Two-phase	2167 MB/s	411 MB/s
256	1 MB	Dynamic Segm.	2491 MB/s	517 MB/s

Tuning parallel I/O performance

- OTPO (Open Tool for Parameter Optimization): optimize the Open MPI parameter space for a particular benchmark and/or application
- Tuning for Latency I/O benchmark on shark/PVFS2
 - Parameters tuned: collective module used, number of aggregators used, cycle buffer size
- 64 different parameter combinations evaluated
- 2 parameter combinations were determined to lead to best performance:
 - *dynamic segm., 20 aggregators, 32 MB cycle buffer size*
 - *static segm. 20 aggregators, 32 MB cycle buffer size*



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sharedfp framework

- Focuses around the management of a shared file pointer
 - Using a separate file and locking
 - Additional process (e.g. mpirun?)
 - Separate files per processes + metadata
 - Shared memory segment
- Collective shared filepointer operations mapped to regular collective I/O operations
- Decision logic based on
 - Location of processes
 - Availability of features (e.g. locking)
 - Hints by the user



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Current status (II)

- Code committed to Open MPI repository in August 2011
- Will be part of the 1.7 release series
- Missing MPI level functionality:
 - Split collective operations (*)
 - Shared file pointer operations: developed in a separate library, currently being integrated with OMPIO (*)
 - Non-blocking individual I/O
 - Atomic access mode



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Part III: Research topics



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OMPIO Optimizations

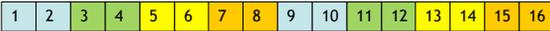
- Automated selection logic for collective I/O modules
- Optimization of collective I/O operations
 - [Development of new communication-optimized collective I/O algorithms \(dynamic segmentation, static segmentation\)](#)
 - [Automated setting of number of aggregators for collective I/O operations](#)
 - [Optimizing process placement based on I/O access pattern](#)
- [Non-blocking collective I/O operations](#)
- [Multi-thread I/O operations](#)



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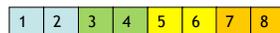
Optimizing communication in collective I/O operations

File layout 

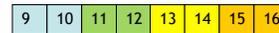
 Process 0  Process 1  Process 2  Process 3

Two-phase I/O with 2 aggregators

Process 0

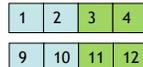


Process 2

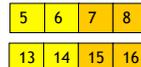


Dynamic segmentation algorithm with 2 aggregators

Process 0



Process 2

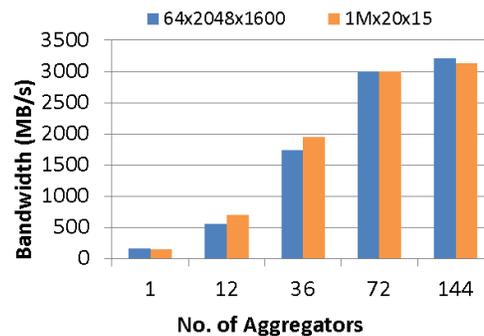


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Automated setting no. of aggregators

- No. of aggregators has enormous influence on performance, e.g.
 - Tile I/O benchmark using two-phase I/O, 144 processes, Lustre file system



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Performance considerations

- Contradicting goals:
 - Generate large consecutive chunks
 - > fewer aggregators
 - Increase throughput
 - > more aggregators
- Setting number of aggregators
 - Fixed number: 1, number of processes, number of nodes, number of I/O servers
 - Tune for a particular platform and application



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Determining the number of aggregators

- 1) Determine the minimum data size k for an individual process which leads to maximum write bandwidth
- 2) Determine initial number of aggregators taking file view and/or process topology into account.
- 3) Refine the number of aggregators based on the overall amount of data written in the collective call

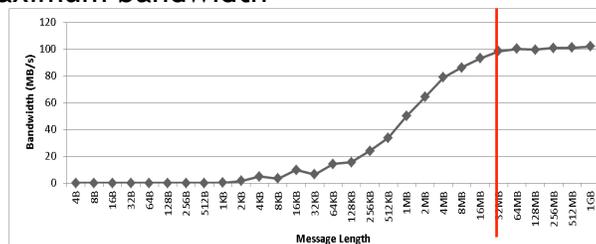


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1. Determining the saturation point

- Loop of individual write operations with increasing data size
 - Avoid caching effects
 - `MPI_File_write()` vs. `POSIX write()`
 - Performed once, e.g. by system administrator
- Saturation point: first element which achieves (close to) maximum bandwidth



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2. Initial assignment of aggregators

- Based on fileview
 - Based on 2-D access pattern
 - 1 aggregator per row of processes
- Based on Cartesian process topology
 - Assumption: process topology related to file access
- Based on hints
 - Not implemented at this time
- Without fileview or Cartesian topology:
 - Every process is an aggregator

Group 1	0	1	2	3
Group 2	4	5	6	7
Group 3	8	9	10	11
Group 4	12	13	14	15



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3. Refinement step

- Based on actual amount of data written across all processes in one collective call
- $k <$ no. of bytes written in group
 - > split group
- $k >$ no. of bytes written in group
 - > merge groups

Group 1	0	1	2	3	Group 2
Group 3	4	5	6	7	Group 4
Group 5	8	9	10	11	Group 6
Group 7	12	13	14	15	Group 8

Group 1	0	1	2	3
	4	5	6	7
Group 2	8	9	10	11
	12	13	14	15



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Discussion of algorithm

- Number of aggregators depends on overall data volume being written
 - Different calls to `MPI_File_write_all` with different data volumes will result in different number of aggregators used
- For fixed problem size, number of aggregators is independent of the number of processes used
- Approach usable for two-phase I/O and some of its variants (e.g. dynamic segmentation)

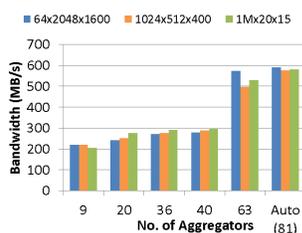


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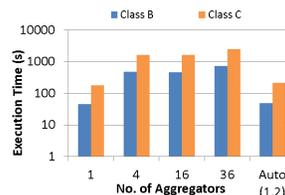


Results

Tile I/O, PVFS2@shark,
81 processes, two-phase I/O



BT I/O, Lustre@deimos,
36 processes, dynamic segmentation



- 134 tests executed in total with 4 different benchmarks
 - 88 tests lead to best or within 10% of optimal performance, 110 within 25% of best performance
- Focusing on two-phase I/O algorithm only:
 - 29 out of 45 test cases outperformed one aggregator per node strategy on average by 41% (default setting by ROMIO)



I/O Access based Process Placement

- Goal: optimized placement of processes to minimize I/O time
- Three required components
 - **Application Matrix:** contains communication volumes between each pair of processes based on the I/O access pattern
 - **Architecture Matrix:** contains communication costs (bandwidth, latency) between each pair of nodes/cores
 - **Mapping Algorithm:** how to map application processes to underlying node architecture such that communication cost are minimized



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Application Matrix

- Goal: predict communication occurring in collective I/O algorithm based on the access pattern of the application
- General case:
 - OMPIO extended to dump the order on how processes access the file
 - Assumption: processes which access neighboring parts of the file will have to communicate with the same aggregators
- Special case:
 - Regular access pattern (e.g. 2D data distribution and process topology)
 - Dynamic segmentation algorithm used for collective I/O
 - Communication occurs only within the outer dimension of the process topology

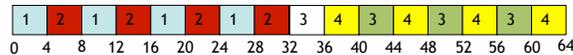


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Application Matrix

- Simple Example : 4 processes with 2x2 tiles, each 4 bytes
- Generic Case: The file layout



- Translates to :

0	7	0	0
7	0	1	0
0	1	0	7
0	0	7	0
- Special Case : Can be represented by topology 2x2 in this case
- Which translates to :

100	100	0	0
100	100	0	0
0	0	100	100
0	0	100	100



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Mapping Algorithms

- Any algorithm from literature could be used
- MPIPP Process Placement Algorithm [1]
 - Randomized algorithm based on Heuristic to exchange processes and calculate gain
 - Generic can support any kind of application and topology matrix
 - Expensive for larger number of processes
- New SetMatch Algorithm for the special case:
 - Create independent sets and matches the sets
 - Very quick even for larger number of processes
 - Greedy approach, and works for specific scenarios
 - Can be generalized by having a clustering algorithm to split sets

[1] Hu Chen, Wenguang Chen, Jian Huang, Bob Robert, and H. Kuhn. 2006. MPIPP: an automatic profile-guided parallel process placement toolset for SMP clusters and multiclusters. In *Proceedings of the 20th annual international conference on Supercomputing (ICS '06)*.



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Preliminary Results

- Crill cluster at the University of Houston
 - Distributed PVFS2 file system using with 16 I/O servers
 - 4x SDR InfiniBand message passing network (2 ports per node)
 - 4x SDR Infiniband (1 port) I/O network
 - 18 nodes, 864 compute cores
- Focusing on collective write operations
- Modified OpenMPI trunk rev. 26077
 - Added a new rmaps component
 - Extensions to OMPIO component to extract fileview information

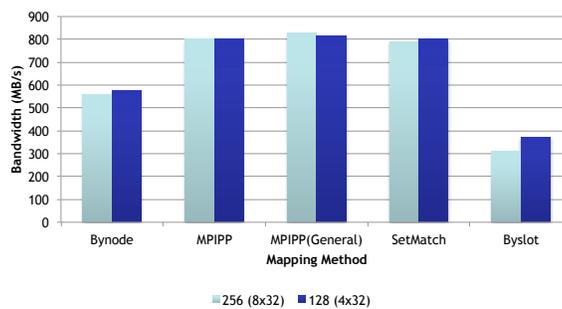


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Tile I/O Results

- Benchmark : Tile I/O
- Tile Size - 1KB
- File size - 128 processes - 75G, 256 processes - 150G

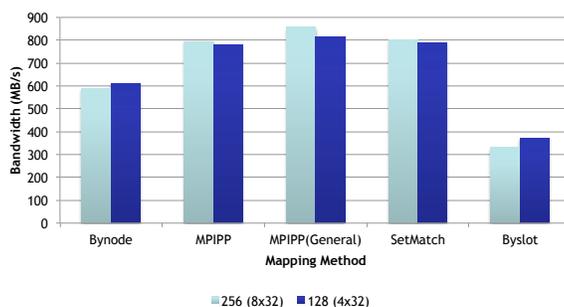


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Tile I/O Results - II

- Benchmark : Tile I/O
- Tile Size - 1MB
- File size - 128 processes - 75G, 256 processes - 150G



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Non-blocking collective operations

- Non-blocking collective Operations
 - Hide communication latency by overlapping
 - Better usage of available bandwidth
 - Avoid detrimental effects of pseudo-synchronization
 - Demonstrated benefits for a number of applications
- Was supposed to be part of the MPI-3 specification
 - Passed 1st vote, failed in 2nd vote

Hoeffler, T., Lumsdaine, A., Rehm, W.: Implementation and Performance Analysis of Non-Blocking Collective Operations for MPI, Supercomputing 2007.

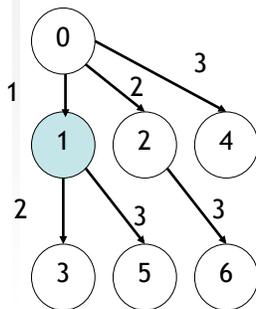


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Overview of LibNBC

- Implements non-blocking versions of all MPI collective operations
- Schedule based design: a process-local schedule of p2p operations is created



Pseudocode for schedule at rank 1:

```
NBC_Sched_rcv(buf, cnt, dt, 0, sched);
NBC_Sched_barr(sched);
NBC_Sched_send(buf, cnt, dt, 3, sched);
NBC_Sched_barr(sched);
NBC_Sched_send(buf, cnt, dt, 5, sched);
```



See <http://www.unixer.de/publications/img/hoefler-hlrs-nbc.pdf> for more details

Overview of LibNBC

- Schedule execution is represented as a state machine
- State and schedule are attached to every request
- Schedules might be cached/reused
- Progress is most important for efficient overlap
 - Progression in `NBC_Test/NBC_Wait`



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Collective I/O operations

- Collective operation for reading/writing data allows to combine data of multiple processes and optimize disk-access
- Most popular algorithm: two-phase I/O
- Algorithm for a collective write operation
 - Step 1:
 - gather data from multiple processes on aggregators
 - Sort data based on the offset in the file
 - Step 2: aggregators write data



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Nonblocking collective I/O operations

```
MPI_File_iread_all (MPI_File file,
                   void *buf, int cnt, MPI_Datatype dt,
                   MPI_Request *request);
```

- Difference to nonblocking collective communication operations:
 - Every process is allowed to provide different amounts of data per collective read/write operation
 - No process has a 'global' view how much data is read/written



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Nonblocking collective I/O operations

- Total amount of data necessary to determine
 - How many cycles are required
 - How much data a process has to contribute in each cycle

⇒ schedule for libNBC can not be constructed in
`MPI_File_iread_all`

- Further consequence:
 - some temporary buffer required internally by the algorithm can not be allocated when posting the operation



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Nonblocking collective I/O operations

- Create a schedule for a non-blocking Allgather(v)
 - Determine the overall amount of data written across all processes
 - Determine the offsets for each data item within each group
- Upon completion:
 - Create a new schedule for the shuffle and I/O steps
 - Schedule can consist of multiple cycles



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Extensions to libNBC

- New internal libNBC operations for:
 - Non-blocking read/write operation
 - Compute operations for sorting and merging entries
 - Buffer management (allocating, freeing buffers)
 - New nonblocking send/rcv primitives with additional level of buffer indirections for dynamically allocated buffers
- Progressing multiple, different types of requests simultaneously



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Caching of schedules

- Very difficult for I/O operations
 - Subsequent calls to `MPI_File_iread_all` will have different offsets into the file
 - Amount of data provided by a process in a cycle depends on the offset in the file
 - Processes allowed to mix individual and collective I/O calls
- ⇒ Not possible to predict offsets of other processes and to reuse a schedule



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Caching of schedules (II)

- When using different files
 - offsets might be the same across multiple function calls, but different file handles will be used
 - Caching typically done on communicator / file handle
- ⇒ Caching across different file handles difficult, but no impossible



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Experimental evaluation

- Crill cluster at the University of Houston
 - Distributed PVFS2 file system using with 16 I/O servers
 - 4x SDR InfiniBand message passing network (2 ports per node)
 - Gigabit Ethernet I/O network
 - 18 nodes, 864 compute cores
- LibNBC integrated with OpenMPI trunk rev. 24640
- Focusing on collective write operations



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Latency I/O tests

- Comparison of blocking and nonblocking versions
 - No overlap
 - Writing 1000 MB per process
 - 32 aggregator processes, 4MB cycle buffer size
 - Average of 3 runs

No. of processes	Blocking Bandwidth [MB/s]	Non-blocking bandwidth [MB/s]
64	703	660
128	574	577



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Latency I/O overlap tests

- Overlapping nonblocking coll. I/O operation with equally expensive compute operation
 - Best case: overall time = max (I/O time, compute time)
- Strong dependence on ability to make progress
 - Best case: time between subsequent calls to `NBC_Test` = time to execute one cycle of coll. I/O

No. of processes	I/O time	Time spent in computation	Overall time
64	85.69 sec	85.69 sec	85.80 sec
128	205.39 sec	205.39 sec	205.91 sec

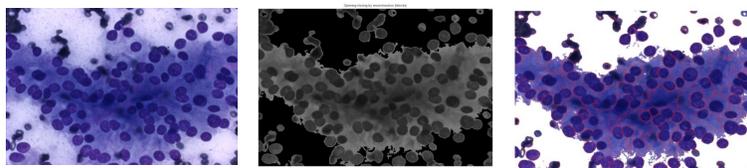


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Parallel Image Processing Application

- Used to assist in diagnosing thyroid cancer
- Based on microscopic images obtained through Fine Needle Aspiration (FNA)
- Slides are large
 - typical image: 25K x 70K pixels, 3-6 Gigabytes/slide
 - multispectral imaging to analyze cytological smears



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Parallel Image Processing Application

- Texture based image segmentation
 - For each Gabor Filter*
 - *Forward FFT of Gabor Filter*
 - *Convolution operation of Filter and Image*
 - *Backward FFT of the convolution result*
 - *Optionally: write result of backward FFT to file*
- FFT operations based on FFTW 2.1.5



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Parallel Image Processing Application

- Code modified to overlap write of iteration i with computations of iteration $i+1$
- Two code versions generated:
 - **NBC**: Additional calls to progress engine added between different code blocks
 - **NBC w/FFTW**: Modified FFTW to insert further calls to progress engine

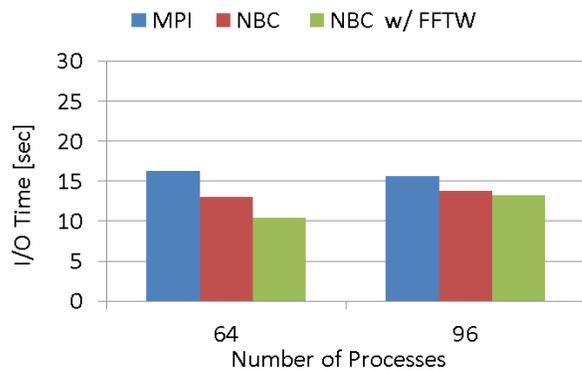


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Application Results (I)

- 8192 x 8192 pixels, 21 spectral channels
- 1.3 GB input data, ~3 GB output data
- 32 aggregators with 4 MB cycle buffer size

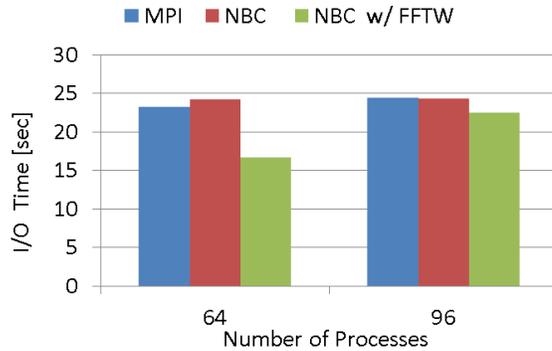


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Application Results (II)

- 12281 x 12281 pixels, 21 spectral channels
- 2.95 GB input data, ~7 GB output data
- 32 aggregators with 4 MB cycle buffer size



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Multi-threaded I/O optimization

- Currently no support for parallel I/O in OpenMP
- Need for threads to be able to read/write to the same file
 - Without locking file handle
 - Without having to write to separate files to obtain higher bandwidth
 - Applicable for all languages supported by OpenMP
- API specification:
 - All routines are library functions (not directives)
 - Routines implemented as collective functions
 - Shared file pointer between threads
 - Support for List I/O Interfaces



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Overview of Interfaces (*write*)

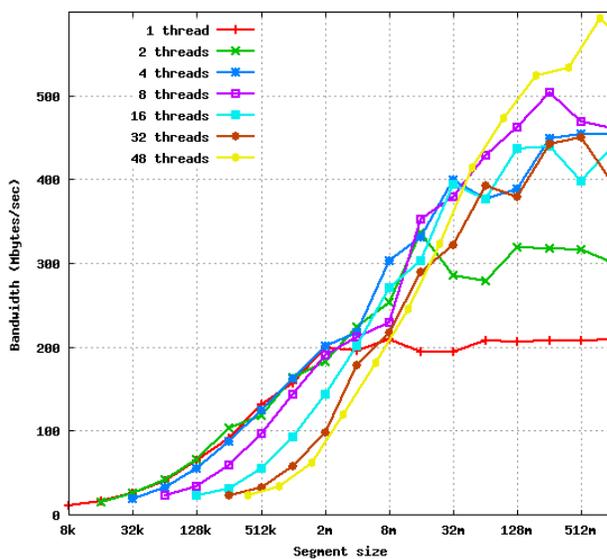
File Manipulation		omp_file_open_all
		omp_file_close_all
Different Arguments	Regular I/O	omp_file_write_all
		omp_file_write_at_all
	List I/O	omp_file_write_list_all
		omp_file_write_list_at_all
Common arguments	Regular I/O	omp_file_write_com_all
		omp_file_write_com_at_all
	List I/O	omp_file_write_com_list_all
		omp_file_write_com_list_at_all



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Results - *omp_file_write_all*



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Performance Results

- OpenMP version of the NAS BT Benchmark
- Extended to include I/O operations

No. of Threads	PVFS2 [sec]	PVFS2-SSD [sec]
1	410	691
2	305	580
4	168	386
8	164	368
16	176	368
32	172	368
48	168	367



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Summary and Conclusions

- I/O is one of the major challenges for current and upcoming high-end systems
- Huge potential for performance improvements
- OMPIO provides a highly modular architecture for parallel I/O
- To improve out-of-the-box performance of I/O libraries
 - Algorithmic developments necessary
 - Handling fat multi-core nodes still a challenge



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