# Lecture 30: Considerations When Using Collective Operations

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### When *not* to use Collective Operations

- Sequences of collective communication can be pipelined for better efficiency
- Example: Process 0 reads data from a file and broadcasts it to all other processes.

```
♦ do i=1,m
if (rank .eq. 0) read *, a
call mpi_bcast(a, n, MPI_INTEGER, 0, comm, ierr)
enddo
```

- Question: How long will this take on p processes?
  - Assume a broadcast takes (s log p + r n) time, and m=p
    - Yes, not (log p) \* (s + rn); the best algorithm is not a distribution tree





#### Broadcast of n Items p Times

- If each takes (s logp + r n) and p
   = m; then the entire loop takes
  - $\bullet$ s \* p log p + p r n
- But there is a way to accomplish this in s p + p r n time!
  - ◆ Log p times as fast if n is small





### Pipeline the Messages

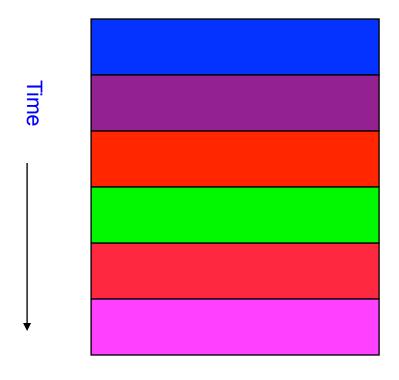
 Process 0 reads data from a file and sends it to the next process. Others forward the data.



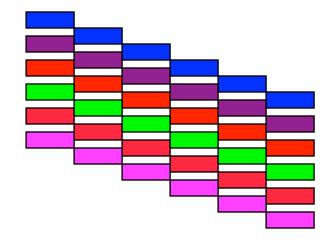
 next = rank+1 unless rank + 1 == size, in which case use MPI\_PROC\_NULL

#### Concurrency Between Steps

Broadcast:



Pipeline



Each broadcast takes less time than pipeline version, but total time is longer



Another example of deferring synchronization

#### Notes on Pipelining Example

- When reading and distributing data from a file, use MPI File read all instead
  - Even more optimizations possible
    - Multiple disk reads
    - Pipeline the individual reads
    - Block transfers
- This algorithm is sometimes called "digital orrery"
  - Circulate particles in n-body problem
  - Even better performance if pipeline never stops
- "Elegance" of collective routines can lead to fine-grain synchronization
  - And hence a performance penalty





#### Thinking about Broadcast

- MPI\_Bcast( buf, 100000, MPI\_DOUBLE, ... );
- Use a tree-based distribution:

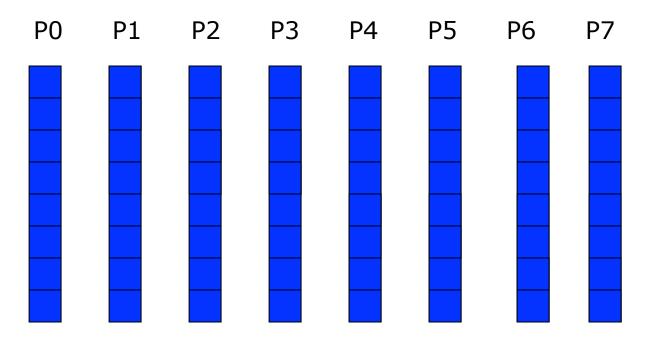


- Use a pipeline: send the message in b byte pieces. This allows each subtree to begin communication after b bytes sent
- Improves total performance:
  - Root process takes same time (asymptotically)
  - Other processes wait less
    - Time to reach leaf is  $b \log p + (n-b)$ , rather than  $n \log p$
- Special hardware and other algorithms can be used ...



#### Make Full Use of the Network

 Implement MPI\_Bcast(buf,n,...) as MPI\_Scatter(buf, n/p,..., buf+rank\*n/p,...) MPI\_Allgather(buf+rank\*n/p, n/p,...,buf,...)







#### Optimal Algorithm Costs

- Optimal cost is O(n) (O(p) terms don't involve n) since scatter moves n data, and allgather also moves only n per process; these can use pipelining to move data as well
  - ◆ Scatter by recursive bisection uses log p steps to move n(p-1)/p data
  - ◆ Scatter by direct send uses p-1 steps to move n(p-1)/p data
  - Recursive doubling allgather uses log p steps to move
    - N/p + 2n/p + 4n/p + ... (p/2)/p = n(p-1)/p
  - Bucket brigade allgather moves
    - N/p (p-1) times or (p-1)n/p
- See, e.g., van de Geijn for more details





#### Implementation Variations

- Implementations of collective operations vary in goals and quality
  - Short messages (minimize separate communication steps)
  - Long messages (pipelining, network topology)
- MPI's general datatype rules make some algorithms more difficult to implement
  - Datatypes can be different on different processes; only the type signature must match



# Using Datatypes in Collective Operations

- Datatypes allow noncontiguous data to be moved (or computed with)
- As for all MPI communications, only the type signature (basic, language defined types) must match
  - Layout in memory can differ on each process





# Example of Datatypes in Collective Operations

- Distribute a matrix from one process to four
  - ◆ Process 0 gets A(0:n/2,0:n/2), Process 1 gets A(n/2+1:n,0:n/2), Process 2 gets A(0:n/2,n/2+1:n), Process 3 gets A(n/2+1:n,n/2+1:n)
- Scatter (One to all, different data to each)
  - ◆ Data at source is not contiguous (n/2 numbers, separated by n/2 numbers)
  - ◆ Use vector type to represent submatrix





### Layout of Matrix in Memory

Process 0

0	8	16	24
1	9	17	25
2	10	18	26
3	11	19	27

32	40	48	56
33	41	49	57
34	42	50	58
35	43	51	59

N = 8 example

Process 2

Process 1

4	12	20	28
5	13	21	29
6	14	22	30
7	15	23	31

36	44	52	60
37	45	53	61
38	46	54	62
39	47	55	63

rocess 3





#### Matrix Datatype

- MPI\_Type\_vector(n/2 per block, n/2 blocks, dist from beginning of one block to next = n, MPI\_DOUBLE\_PRECISION,&subarray\_type)
- Can use this to send
  - ◆ Do j=0,1 Do i=0,1 call MPI\_Send( a(1+i\*n/2, 1+j\*n/2), 1, subarray\_type, ... )
  - ◆ Note sending ONE type contain multiple basic elements
  - Pass the (address of the) first element to be sent to MPI\_Send
  - This looks like an MPI\_Scatter, but with interleaved data





#### Scatter with Datatypes

- Scatter is like
  - ◆ Do i=0,p-1 call mpi\_send(a(1+i\*extent(datatype)),....)
    - "1+" is from 1-origin indexing in Fortran
  - Extent is the distance from the beginning of the first to the end of the last data element
  - ♦ For our subarray\_type, it is ((n/2-1)n+n/2) \* extent(double)
  - "extent(double)" is simply the number of bytes in DOUBLE PRECISION item (often 8)
    - In Fortran, you can use MPI\_Type\_size( MPI\_DOUBLE\_PRECISION, extent, ierr)
    - Or MPI\_SIZEOF(a) (with the MPI or MPI\_F08 module)
    - Or storage size(1.0do)/8 (in Fortran 2008)







# If Only We Could Change the Extent of subarray\_type...

- To make the communication work with Scattery, set Extent of each datatype to n/2
  - Size of contiguous block all are built from
- Use Scattery (independent multiples of extent)
- Location (beginning location) of blocks
  - Process 0: 0 \* 4 (doubles)
  - ◆ Process 1: 1 \* 4 (doubles)
  - Process 2: 8 \* 4 (doubles)
  - Process 3: 9 \* 4 (doubles)
- How can we change the extent of a datatype?





### Changing the Extent

- MPI allows you to change the extent of a datatype with MPI\_Type\_create\_resized
- In our case (in C),
- MPI\_Type\_create\_resized( subarray\_type, 0, (n/2)\*sizeof(double), &newtype)
  - ◆ Sets the lower bound to zero (almost always the right thing) and the extent to n/ 2 doubles.





#### Scattering A Matrix

```
    sdisplace(1) = 0
        sdisplace(2) = 1
        sdisplace(3) = n
        sdisplace(4) = n + 1
        scounts(1,2,3,4)=1
        call MPI_Scatterv(a, scounts, sdispls, newtype, & alocal, n*n/4, MPI_DOUBLE_PRECISION, & 0, comm, ierr)
```

- ◆ Note that process 0 sends 1 item of newtype but all processes receive n²/4 double precision elements
- Test yourself: Work this out and convince yourself that it is correct





### Dense Matrix Vector Multiply

- Let the matrix be partitioned across processes by columns, and the vector by corresponding rows.
  - ◆ If process i has columns M:N of the matrix, it also has elements M:N of the vectors
  - ◆ Simple partition (process 0 has the first block of columns, process 1 the second block, etc.)
  - process i has columns col(i):col(i+1)-1
- Problem: Compute the matrix-vector product with the distributed data structures
  - Send/receive requires intermediate buffers
  - Collective solution





#### Using MPI\_Reduce\_scatter

 Each process needs to accumulate a contribution from every process to its part of the result vector





### Meaning of Reduce Scatter

- Reduce\_scatter
  - Scatters contributions from all processes to all others
  - Combines (reduces) incoming contributions into a single buffer
  - MPI\_Reduce\_scatter\_block like MPI\_Reduce\_scatter, but with the same size block on all processes
- Reduce\_scatter also be used for distributed inmemory checkpoint with error correction
  - See SCR <a href="https://computation.llnl.gov/project/scr/">https://computation.llnl.gov/project/scr/</a>
  - ◆ Providing Efficient I/O Redundancy in MPI Environments, Gropp, Ross, Miller, EuroPVM/MPI 2004,



http://link.springer.com/chapter/ 10.1007/978-3-540-30218-6\_17



#### Some Performance Issues

- MPI Collectives must handle the general case
- Implementations usually optimize for each collective operation separately
  - Assumption is make each individual collective as fast as possible, not the overall application
  - ◆ A Study of Process Arrival Patterns for MPI Collective Operations, Faraj, Patarasuk, Yuan, IJ Parallel Programming, 36:6 2008
    <a href="http://link.springer.com/article/10.1007%2Fs10766-008-0070-9">http://link.springer.com/article/10.1007%2Fs10766-008-0070-9</a>
- Implementations sensitive to progress (availability of CPU to advance communication)



- Particularly important for nonblocking collectives
- Nonblocking doesn't ensure concurrent execution

