Distributed Computation via MPI

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There are two common parallelization models in parallel programming

- Master / Worker (like what we have seen in parallel's parLapply)
- Single Program Multiple Data (SPMD)

SPMD (single program, multiple data) is a technique employed to achieve parallelism; it is a subcategory of MIMD. Tasks are split up and run simultaneously on multiple processors with different input in order to obtain results faster.

Advantages of SPMD over Master/Worker - It is very close to the serial code. i.e. SPMD is easy to modify from serial. - It is much shorter than the original Master/Worker version. i.e. SPMD is traceable for debugging. - It makes the master as one of workers. i.e. SPMD fully utilizes resources. - It is easy to automatically process large numbers of independent jobs. i.e. SPMD can parallelize by jobs.

MPI

MPI (Message Passing Interface) is one of the most standard parallel computing architectures. In R, there are two MPI packages Rmpi and pbdMPI.

- Rmpi has been here for a long time, so it is more popular to Rmpi given its age. However it is not actively maintained now which means that Rmpi is not working probably with new versions of MPI (for example, >= openmpi 2.0).
 - Rmpi could be used to execuate both master/worker type of jobs or SPMD type of jobs but there is a bug in Rmpi with makes master/worker mode fails in our cluster.
- pbdMPI is a late comer (which means that it is actively maintained) and it claims that it's fater than Rmpi

SPMD

The general process in SPMD programs goes something like this:

- initialize communicators
- have each process read in its portion of the data
- perform computations
- communicate results
- shut down the communicators

Our cluster

Our cluster uses SLURM to manage resources. There are two common resource management systems, PBS and SLURM. It's more difficult to adapt to the other system once you know how a system works.

- first ssh into your account on peloton.cse.ucdavis.edu
- upload your R script to the server (either by a GUI software such as Cyber Duck/ WinScp or using scp)
- Use the following to download all the demo codes,

svn export https://github.com/UCDavis-STA-141C-Winter-2020/sta141c-lectures/branches/master/02-11/demo/

Important! - Do not run R in the head node - Be considerate, do not specific more than enough cores. Our class has 128 cores shared between 50 students. Use sinfo to query the current status of the cluster and squeue to check the running jobs in the cluster.

Basic Communicator Wrangling

Managing a Communicator: Create and destroy communicators - init() Initialize communicator - finalize() shut down communicator(s) Rank query: Determine the processor's position in the communicator - comm.rank() "who am I?"" - comm.size() "how many of us are there?"" - Print: printing with control over which processor prints. - comm.print(x, ...) - comm.cat(x, ...)

To run the script O1-pbdmpi.R

```
# make site packages such as pdbMPI available
module load R
# use 3 cores and 1 minute limit
# high2 is a partition reserved for our class
srun -p high2 -n 3 -t 1 Rscript 01-pbdmpi.R
# use 3 cores on 2 nodes and 1 minute limit
srun -p high2 -N 2 -n 3 -t 1 Rscript 01-pbdmpi.R
```

You could press ctrl+c anytime to interrupt the job.

Communication functions

- bcast A Rank Broadcasts an Object to Every Rank
- scatter A Rank Scatter Objects to Every Rank
- gather/allgather A Rank / All Ranks Gather(s) Objects from Every Rank
- reduce/allreduce A Rank / All Ranks Receive(s) a Reduction of Objects from Every Rank
- send/recv A Rank Sends / Receives (blockingly) an Object to / from the Other Rank
- isend A Rank Sends (non-blockingly) an Object to the Other Rank

Parallel Lapply Functions

- pbdLapply analogue of lapply()
- pbdSapply analogue of sapply()

There are three modes: mw, spmd and dist.

Reference:

- \bullet Data Science with R and pbdR at ORNL https://www.olcf.ornl.gov/wp-content/uploads/2018/07/pbdR_part2.pdf
- A Forgotten MPI Tool For R http://thinkdatascience.com/pbdr-a-forgotten-mpi-tool-for-r/