PHYS405

Advanced Computational Physics Parallel Computing

Assignment # 2Due: Friday, October 21, 2016

Purpose: Learn to write and run a simple MPI code

Note: Please identify all your work.

Write an MPI code, submit.c, to execute a Linux command on all the nodes of a virtual machine. The command to execute will be extracted from the command line arguments.

For instance,

mpiexec -np 4 ./submit rm /tmp/a

will remove the file \mathbf{a} from the local directory $/\mathbf{tmp}$ on all the nodes on which the code runs.