

High Performance Optimizations and Dynamic Load Balancing for Computational Aerodynamics Solvers

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1 Initial Requirements

1. Use the Ubuntu operating system and your PC must have at least 4 cores.
2. You must install the OPEN MPI latest version using either of the following two ways.
 - `sudo apt-get install openmpi-bin openmpi-common openssh-client openssh-server libopenmpi1.3 libopenmpi-dev libopenmpi-dev`

- Download the open-mpi tar file from <http://www.open-mpi.org/software/ompi>

- Decompress the downloaded file (should be called something like `openmpi-x.x.x.tar.xxx`, changing `x.x.x` for the version downloaded):

```
tar -xvf openmpi-*
```

- Go into the new folder created from the decompress.

```
cd openmpi-*
```

- It is necessary to add on the prefix the installation directory we want to use for OpenMPI. The normal thing to do would be to select the next directory `"/home/user/.openmpi"`.

```
./configure --prefix="/home/$USER/.openmpi"
```

- Install the MPI

```
make
sudo make install
```

- All that is left to do is to include the path to our path environment the path “installation_directory/bin” and to the library environment variable “installation_directory/lib/”. If your system uses bash you’ll have to use the command export.

```
export PATH="PATH : /home/USER/.openmpi/bin"
export LD_LIBRARY_PATH="LD_LIBRARY_PATH : /home/USER/.openmpi/lib/"
```

3. Next you must have metis software package installed on your system.

- Download the metis tar file from
<http://www.gilith.com/software/metis/download.html>
- Extract it with the command

```
tar xvzf metis.tar.gz
```

- Change to directory metis

```
cd metis
```

- Building METIS requires CMake 2.8, found at <http://www.cmake.org/>, as well as GNU make. Assuming CMake and GNU make are installed, two commands should suffice to build metis:

```
make
make config
```

- Install the metis

```
make install
```

- Change into the directory as

```
cd /build/Linux-x86_64/programs/
```

- In this folder you can see a file 'gpmetis' which is executable and you can use this program to partition the graph using the following command

```
./gpmetis filename.graph count
```

4. You must have a gcc compiler and g++ compiler in your system you can install it using

```
sudo apt-get install gcc
sudo apt-get install g++
```

2 Running the system

1. Clone into the following repository

```
git clone https://github.com/robinl3680/Final-Year-Project.git
```

2. Change into the directory

```
cd Final-Year-Project/Progress/MeshToGraph
```

3. Compile and Execute the Mesh to Graph conversion algorithm as follows.

```
g++ MeshToGraph.cpp -o MtoG  
./MtoG input_file_name.su2 output_file_name.graph output_faces.txt
```

4. Go to the directory 'MPI'

```
cd ../MPI/
```

5. In order to execute the different communication algorithms individually do the following.

- Centralized communication approach

```
mpicxx Centralized.cpp -o Centralized  
mpirun -np cores_count+1 Centralized input.graph.part.count input.graph
```

- Peer to Peer communication approach

```
mpicxx PeerToPeer.cpp -o Peer  
mpirun -np cores_count Peer input.graph.part.count input.graph
```

- Greedy communication approach

```
mpicxx Greedy.cpp -o Greedy  
mpirun -np cores_count Greedy input.graph.part.count input.graph
```

- Another way to run the three in once use following shell script and you can edit the input files mentioned there.

```
sh testbash.sh
```

3 Running within a Cluster

1. Connect the different systems using a wired or wireless LAN.
2. Make sure that all system contain same version of the 'MPI'. To check that.

```
mpirun -version
```

3. Using the ping command check the physical connectivity between systems.

```
ping ip_address
```

4. Make sure that each system has the required programs to be executed are in the same path.
5. Enable password less login between the systems to avoid the authentication issues.

```
ssh-keygen -t rsa  
ssh-copy-id remote_username@server_ip_address
```

6. Check the \$PATH variable to ensure they are same.

```
echo $PATH
```

7. If they are not same make it same by copying one of them and replacing it with others using

```
export $PATH=new_path
```

8. Create a host file that contain the ip addresses of the connected systems.

```
nano host_file  
ip_address1  
ip_address2  
...
```

9. Now execute the different communication algorithms by specifying the host file. An example is.

```
mpicxx Greedy.cpp -o Greedy  
mpirun -np cores_count -hostfile host_file_name Greedy input.graph input.graph.part.count
```