

## **EXE – PowerFlow – Execution Guidelines for running applications in aggregated environment using ScaleMP’s vSMP Foundation**

### ***Overview***

PowerFlow’s simulation phase is a multi-process application that uses MPI for inter-process communication. HP-MPI has been set as the default MPI for the PowerFlow application. In addition upcoming versions of PowerFlow would support Intel-MPI/MPICH2 as well.

Guidelines for running PowerFlow with Intel-MPI / MPICH2 would be available as soon as the new version of PowerFlow is released.

### ***Running PowerFlow with ramfs***

Since PowerFlow jobs perform a lot of scratch IO, running with a ramfs may yield shorter runtimes. In order to run with ramfs, perform the following:

- 1) Make sure that /ramfs is mounted, and if not mount it as follows:

```
sudo mkdir /ramfs

sudo mount /dev/ram -t ramfs /ramfs -o noatime

sudo chmod 777 /ramfs
```

- 2) Copy the input into /ramfs and run from there, so the '\*.lgi' file would be created in /ramfs

### ***Separating between the discretizer and Simulation***

The discretizer part is not intended to run on a large number of cores; while the simulation part is. In order to run it optimally, one should run the discretizer phase separately with up to 8 cores, and bound this part of the job to a single board, by using:

```
numactl -cpunodebind=2,3 [run script]
```

The above example would bind the process to the 2<sup>nd</sup> board.

After the discretizer phase is complete, run the simulation part using MPI as described below.

### ***Number of Ranks***

As PowerFlow always creates one additional control process; when the target is to launch it using N cores, please specify 'N-1' as the number of MPI ranks to use.

## Running PowerFlow with HP-MPI

HP-MPI has a built-in mechanism for assigning MPI processes to specific CPUs. Process placement is controlled by environment variables named **MPI\_BIND\_MAP** and **MPIRUN\_OPTIONS**. When these variables are not set, process placement will not be performed.

### Environment variables – HP-MPI

If you are running Fluent with HP-MPI, you should set the following environment variables prior to running Fluent to yield the optimal performance:

```
export MPI_BIND_MAP=0,1,2,3,4,5,6,7 (For example)

export MPIRUN_OPTIONS="-cpu_bind=map_cpu,v"

export HPMP_FRAGSIZE=131072

export MPI_SHMEMCNTL=16,24000000,4000000
```

MPI\_BIND\_MAP specifies a list of CPUs to which MPI ranks will be bound. You should replace the list above with a list of integers, zero to #cpus-1.

For more information on HP-MPI CPU affinity settings, refer to the HP-MPI user's guide available from ["http://docs.hp.com/en/B6060-96022/B6060-96022.pdf"](http://docs.hp.com/en/B6060-96022/B6060-96022.pdf).

### Sample script for HP-MPI

```
#!/bin/sh

export PATH=/opt/powerflow/4.2c/bin:$PATH
export LM_LICENSE_FILE=27007@<hostname>
export NCPUS=`grep -i processor /proc/cpuinfo | wc -l`
export MPI_BIND_MAP=`seq -f "%g," 0 $((NCPUS-1)) | tr -d "\n" | sed -e 's/,,$//`
export MPIRUN_OPTIONS="-cpu_bind=map_cpu,v"
export HPMP_FRAGSIZE=131072
export MPI_SHMEMCNTL=16,24000000,4000000

# Always, use Nproc-1 processes as PowerFlow creates one additional control process.
np=15

taskset 0xff exaqsab -nprocs $np external_case_2.cdi
```

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***Running PowerFlow with MPICH2 tuned for vSMP***

Would be available as soon as the new version of PowerFlow is released.