

High Performance Computing with Python Execution Instructions

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Execution Instructions

To execute locally:

1. Create an anaconda environment as follows:

```
conda create --name hpc python=3.8 -y
conda activate hpc
pip install -U pip
pip install -r requirements.txt
```

- 2. Navigate into src directory.
- 3. Execute the sliding_lid_parallel.py using mpirun and the following command line arguments:
 - Use the mpirun -n command line argument to specify the number of cores.
 - Use the -g command line argument of sliding_lid_parallel. py to specify the lattice's x- and y-dimensions.
 - Use the -d command line argument of sliding_lid_parallel.
 py to specify the desired domain decomposition in the x- and y-dimensions.
 - Use the -o command line argument of sliding_lid_parallel. py to specify the collision frequency.
 - Use the -v command line argument of sliding_lid_parallel. py to specify to specify the lid velocity.

Example Execution

Execute the sliding lid simulation in parallel on a 300×300 grid with collision frequency equal to 1.7, lid velocity equal to 0.1 for 10000 steps. Use 4 cores and 2×2 discretization.

```
mpirun -n 4 python sliding_lid_parallel.py -g 300 300 -d 2 2 -o 1.7 -v 0.1
```

To get further information on the command line arguments available for sliding_lid_parallel.py execute:

```
python sliding_lid_parallel.py -h
```

The execution will create two npy files for the components of velocity. Execute the following script to visualize the end result

```
python visualize_velocity_field_ilias.py
```

For additionally information on the parallel execution as well as the execution of the rest of the experiments please advise the report's Github repository https://github.com/theodorju/fr-hpcpy-pub.