## **README**

Docariation

## Description

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The programme developed is responsible for generating appropriate initial conditions and solving a two-dimensional smoothed particle hydrodynamic (SPH) formulation of the Navier-Stokes equations. The files that participate in that process are:

- main-SPH.cpp
- class.cpp
- class.h
- Makefile

The running process will result in two files:

- One for the Energies at each time step
- One for the final positions of the particles

Compiling

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In order for the user to compile the programme, the command "make" has to be typed in the command line. The compiling process will result in an executable file called "SPH-SOLVER".

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Proper execution

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Upon executing "SPH-SOLVER", the user should insert four command line arguments:

- One for the initial condition (--ic-dam-break, --ic-block-drop, --ic-droplet, --ic-one-particle, --ic-two-particles, --ic-three-particles, --ic-four-particles)
- One for time step to use (--dt)
- One for the total integration time (--T)
- One for the radius of influence of each particle (--h)

Example of running the executable for dam-break, timestep of 0.0001, total integration time of 3 seconds and a radius of influence 0.01:

mpiexec -np 2 ./SPH-SOLVER --ic-dam-break --T=3 --dt=0.0001 --h=0.01

The code has a specific number of particles in it for each initial condition. Supposing the users would like to change the number of particles they should be very careful, since the dam-break and droplet conditions can handle only numbers of particles that are a power of 2. Also, the number of particles in block-drop initial condition is carefully chosen, so that dx=dy. These practices were used to make sure that a uniform distribution is achieved across the whole grid.

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Possible Issues

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Generally, the code responds well to the use of many processors. However, for unidentified reasons some cases provided errors, while on a different day they may have executed the same simulation without any problems. It is advised to use 5 or less processors to avoid problems