

# Event driven simulation of a 2-dimensional gas

**Sondre Duna Lundemo<sup>†</sup>**

Department of Physics, Norwegian University of Science and Technology, Trondheim  
Norway

TFY4235 - Computational physics

*(Last updated on January 29, 2021)*

An event driven approach is used to simulate a gas of hard discs in two dimensions. The framework built is used to study statistical properties of the gas at and towards equilibrium. The numerical results are compared with well known results from classical statistical mechanics.

# Contents

<b>1</b>	<b>Introduction and background</b>	<b>2</b>
<b>2</b>	<b>Overview of code</b>	<b>2</b>
2.1	Remarks on performance . . . . .	3
<b>3</b>	<b>Results</b>	<b>3</b>
3.1	Speed distribution at equilibrium . . . . .	3
3.2	Mixture of two gases . . . . .	4
3.3	Mixing in the presence of energy dissipation . . . . .	4
3.4	Crater formation from projectile impact . . . . .	4

# 1 Introduction and background

An event driven approach is used to simulate a gas of hard discs in two dimensions. The framework built is used to study statistical properties of the gas at and towards equilibrium. The numerical results are compared with well known results from classical statistical mechanics.

The central algorithm for event-driven simulations is the following [1]:

```

Set velocities and positions of all particles in the gas;
Choose a stop criterion ;
for each particle in gas do
    Calculate if and when the particle will collide with all the other particles and
    the walls ;
    Store all the collision times;
end
while not reached stop criterion do
    Identify the earliest collision ;
    if collision is valid then
        Move all particles in straight lines until the earliset collision ;
        for each particle involved in collision do
            Calculate if and when the particle will collide with all the other
            particles and the walls ;
            Store all the collision times;
        end
    else
        discard collsion ;
    end
end

```

**Algorithm 1:** Event driven simulation of a gas.

In the above algorithm, a collision is *valid* if the particle(s) involved in the collision has *not* collided since the time the collision time was stored.

# 2 Overview of code

In this section we briefly present how the algorithm is implemented and which considerations are put into the choices of data-structures.

## 2.1 Remarks on performance

# 3 Results

## 3.1 Speed distribution at equilibrium

The first test is devoted to investigating the distribution of the velocities of the particles after sufficiently long time. From statistical mechanics, we know that the velocities of a 2-dimensional gas at equilibrium will distribute according to Maxwell's velocity distribution:

$$p(v) = \frac{mv}{k_B T} \exp\left(-\frac{mv^2}{2k_B T}\right),$$

where  $k_B$  is Boltzmann's constant,  $T$  the absolute temperature, and  $m$  the mass of the particles. It is essential that the all the particles have the same mass  $m$ , as will become apparent in the following section [LINK].

We initialise an ensemble of particles with velocities  $\mathbf{v} = [v_0 \cos \theta, v_0 \sin \theta]$  with  $\theta \sim \mathcal{U}[0, 2\pi]$ . The initial distribution of the velocities is therefore  $C\delta(v - v_0)$ . After simulating sufficiently long, the distribution is as shown in figure ??.

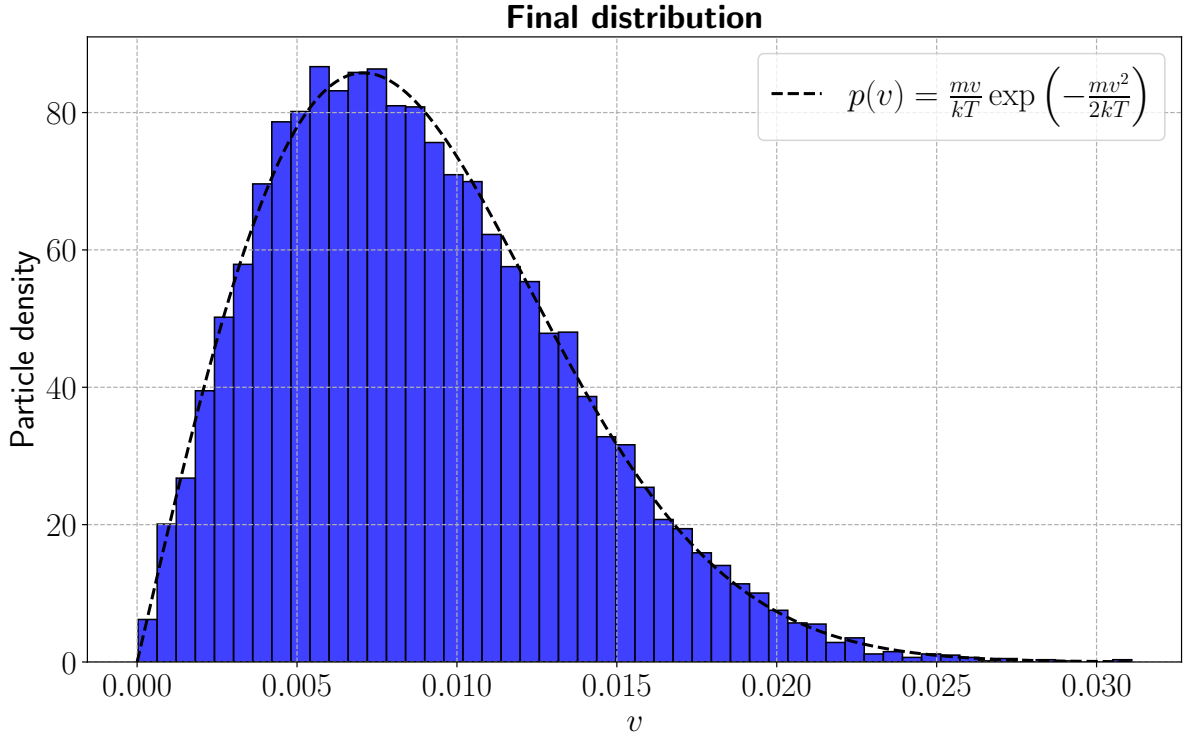


Figure 1: Distribution of velocities in an ensemble of 10000 particles after 200 s.

### 3.2 Mixture of two gases

### 3.3 Mixing in the presence of energy dissipation

### 3.4 Crater formation from projectile impact

## References

- [1] Aleksandar Donev, Salvatore Torquato, and Frank H. Stillinger. Neighbor list collision-driven molecular dynamics simulation for nonspherical hard particles. i. algorithmic details. *Journal of Computational Physics*, 202(2):737 – 764, 2005.