

Advanced Computer Graphics

Homework 6

Particle System Part 2

Group 8
Damien Doy - 244206
Jason Racine - 244270
Alexandre Veuthey - 224295

Computation of the triangle area forces

In order to find the force at each vertice of the triangle we have to find the following partial derivative (as stated in the lecture):

$$F_i = -\frac{\partial EA(x_0, x_1, x_2)}{\partial x_i} \text{ where } EA(x_0, x_1, x_2) = 0.5k_A(\text{area}(x_0, x_1, x_2) - A)^2$$

and

$$\text{area}(x_0, x_1, x_2) = 0.5((x_1.x - x_0.x)(x_2.y - x_0.y) - (x_2.x - x_0.x)(x_1.y - x_0.y))$$

By solving these equations, we get:

$$F_i = -\frac{\partial EA(x_0, x_1, x_2)}{\partial x_i} = k_A(\text{area}(x_0, x_1, x_2) - A) \frac{\partial \text{area}(x_0, x_1, x_2)}{\partial x_i}$$

And by solving the partial derivative we get the following forces for each particles:

$$F_0 = -0.5dEA(x_{y1} - x_{y2}, x_{x2} - x_{x1})$$

$$F_1 = -0.5dEA(x_{y2} - x_{y0}, x_{x0} - x_{x2})$$

$$F_2 = -0.5dEA(x_{y0} - x_{y1}, x_{x1} - x_{x0})$$

$$\text{where } dEA = k_A(\text{area}(x_0, x_1, x_2) - A)$$

Computation of the Jacobian

Gravity

Let's begin with the gravitation force, which is defined as

$$F_G = \begin{pmatrix} 0 \\ -Gm \end{pmatrix}$$

Since the force of gravity does not depend on the position of the particle, the 2x2 Jacobian of this force by the particle position is a null matrix and doesn't have to be added to the solver

Centre force

The center force is defined as:

$$F_C = \begin{pmatrix} -Cx \\ -Cy \end{pmatrix}$$

And thus, the jacobian of this force is:

$$J_C = \begin{pmatrix} -C & 0 \\ 0 & -C \end{pmatrix}$$

Force-based collisions

The force based collision is applied to the particle only when

$xn_x + yn_y + n_z - p_{radius} < 0$ where x and y are the position of the particle, and n_x , n_y and n_z are the parameters of the line equation. Then the force is:

$$F_{Col} = \begin{pmatrix} -n_x(xn_x + yn_y + n_z - p_{radius})k_{col} \\ -n_y(xn_x + yn_y + n_z - p_{radius})k_{col} \end{pmatrix} = \begin{pmatrix} -k_{col}(xn_x^2 + yn_xn_y + n_xn_z - p_{radius}n_x) \\ -k_{col}(xn_xn_y + yn_y^2 + n_yn_z - p_{radius}n_y) \end{pmatrix}$$

Then the jacobian is

$$J_{Col} = \begin{pmatrix} -k_{col}n_x^2 & -k_{col}n_xn_y \\ -k_{col}n_xn_y & -k_{col}n_y^2 \end{pmatrix}$$

Damped-spring forces

Forces are:

$$F_{spring0} = - \begin{pmatrix} (k_s(||p_0 - p_1|| - L) + k_d \frac{(x_0 - x_1) \cdot (x_0 - x_1)}{||p_0 - p_1||}) \frac{x_0 - x_1}{||p_0 - p_1||} \\ (k_s(||p_0 - p_1|| - L) + k_d \frac{(y_0 - y_1) \cdot (y_0 - y_1)}{||p_0 - p_1||}) \frac{y_0 - y_1}{||p_0 - p_1||} \end{pmatrix}$$

This force can be cumbersome to calculate, so we separated it into 3 components:

$$Stiffness_x = k_s(||p_0 - p_1|| - L)$$

$$Damping_x = k_d \frac{(x_0 - x_1) \cdot (x_0 - x_1)}{||p_0 - p_1||}$$

$$Vector_x = \frac{y_0 - y_1}{||p_0 - p_1||}$$

Here are the three component for forces in x only, but the same applies to the forces in y.

By taking the partial derivative for all four variables, for the three component, we can then calculate the jacobian for the damped-spring force in the form:

$$\frac{\partial}{\partial x_0} F_{spring0x} = \frac{\partial}{\partial x_0} - (Stiffness + Damping)Vector$$

This makes the jacobian much easier to calculate and less susceptible to errors.

The result jacobian is a 4X4 matrix as there are two particles acting on two forces:

$$\frac{\partial F_{spring0}}{\partial p_0}, \frac{\partial F_{spring1}}{\partial p_0}, \frac{\partial F_{spring0}}{\partial p_1} \text{ and } \frac{\partial F_{spring1}}{\partial p_1}$$

This calculation is done in the code in the function `compute_jacobians` and the same principle has been applied to the damped spring forces of the mouse click except here,

Area preservation forces

One of the force is defined as:

$$F_{area0} = -0.5dEA \begin{pmatrix} p_{y1} - p_{y2} \\ p_{x2} - p_{x1} \end{pmatrix}$$

The other forces are defined at the first chapter where we defined how we calculated the forces. Since we have 3 forces acting on 3 particles, this means that we have a 3X3 jacobian matrix. The way we calculated it is the same as the damped string forces. The calculation can get messy and won't be described here.

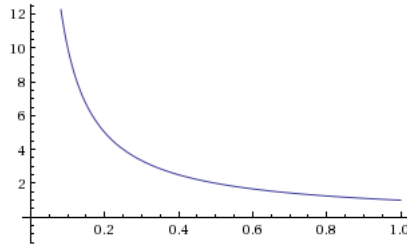
Computation of the equilibrium force

In order to distribute the particles on the screen, we added a force that acts on every particles. They repulses themselves like same polarity charges in electromagnetism.

To avoid a unstable system and to reduce the computation time, we limited this force to a certain radius around each particle. This radius is equals to

$radius = C(\frac{1}{dist} - 1) * normal_{p_0, p_1}$ where $dist \in [0, 1]$ is the distance between two particles normalized. C is a constant that makes the equilibrium force stronger or weaker.

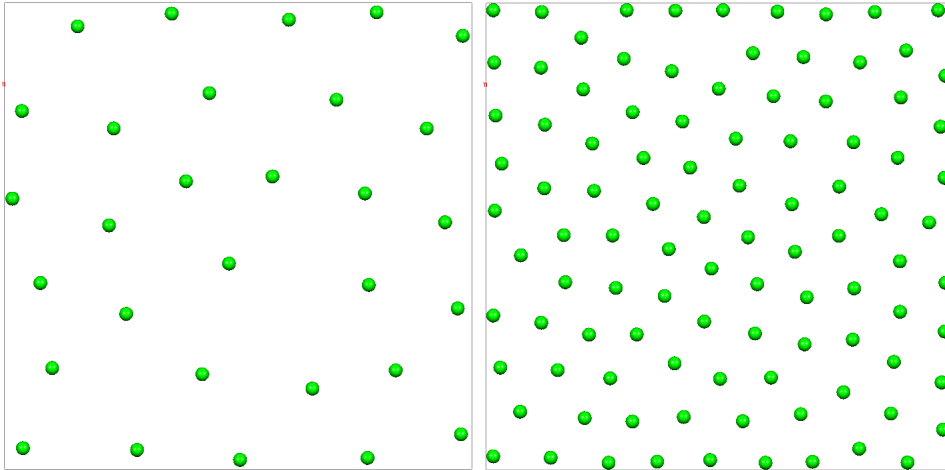
Here is a graph of this function:

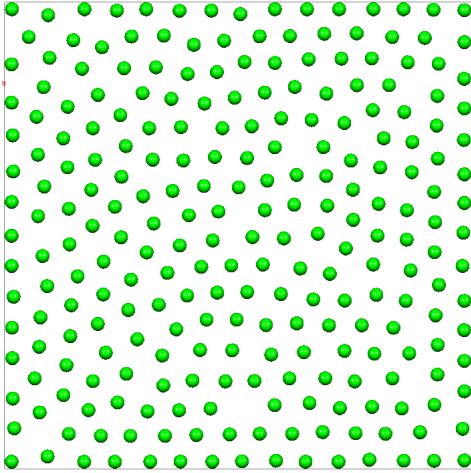


The advantage of a strong force, is that particles will distribute uniformly faster, but a great force will make the system go unstable.

A weak force can also be advantageous, particles will move slower to their positions but there is less risk of particles flying off.

Here we tested the equilibrium force on 30, 100, 250 particles, with verlet integration and impulse based forces:





Comparison of the integration methods

The verlet integration is the most stable of the integrations methods in this lab.

The euler integration method is really unstable and particles will rarely find a rest point (they will bounce around or jiggle when there are springs). The system can go unstable very easily with springs. A lower timestep will solve this problem at the cost of computation time.

The midpoint integration method have the same problems as euler but the instabilities are reduced for the same time stamp.

In the case of euler and midpoint integration, all inaccuracies and the instability of the system can be solved by reducing the timestamp.

The verlet integration is the most stable. For the same timestamp as used for euler and midpoint, particles will find a rest point at some point and the scene with the octagon will stay stable, even if violent forces are acting on the shape. The verlet integration can be even used with larger timestamps, the system will stay stable at $dt=0.002$ and will show some problems at $dt=0.004$.

In the case of implicit euler, the same problems as explicit euler appear. That is, the system go unstable when the timestamp is too high. A lower time stamp ($t=0.00025$) solve the problem.