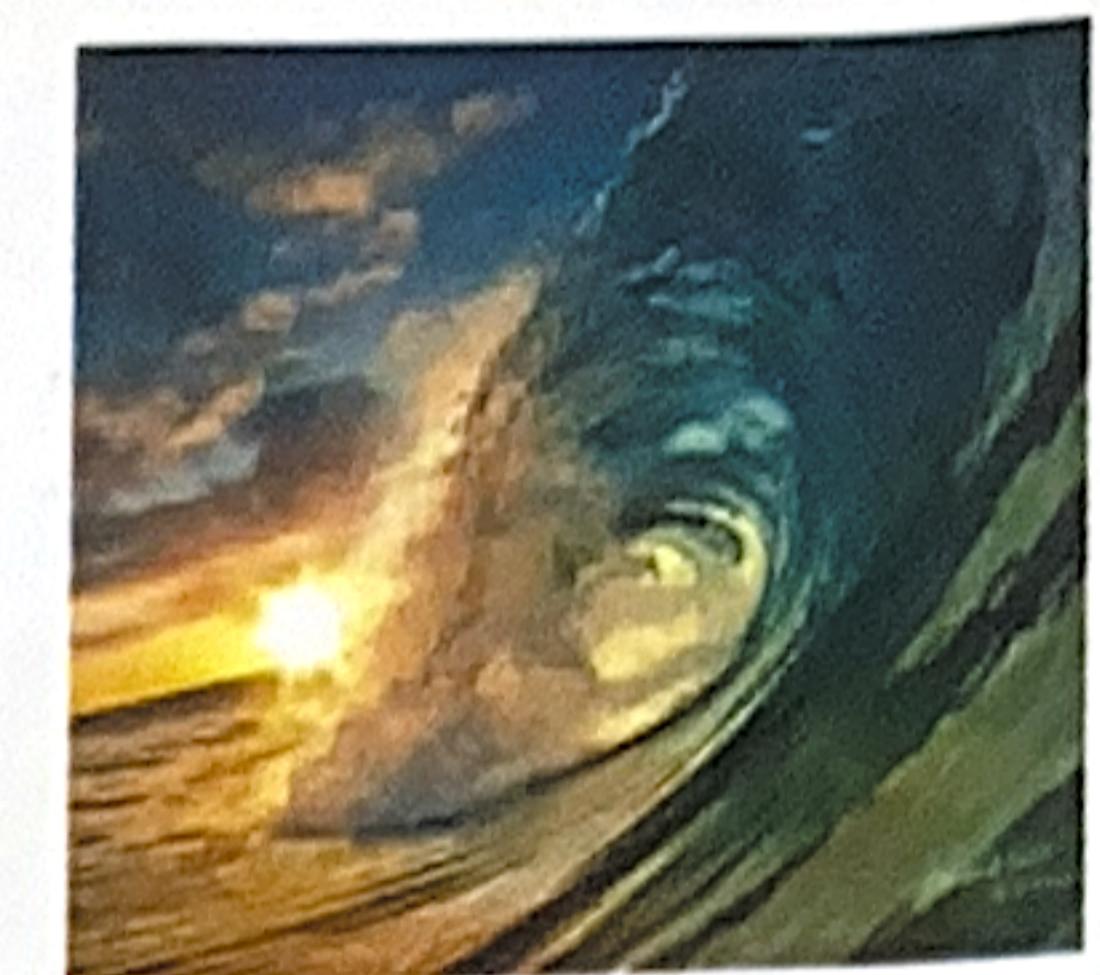
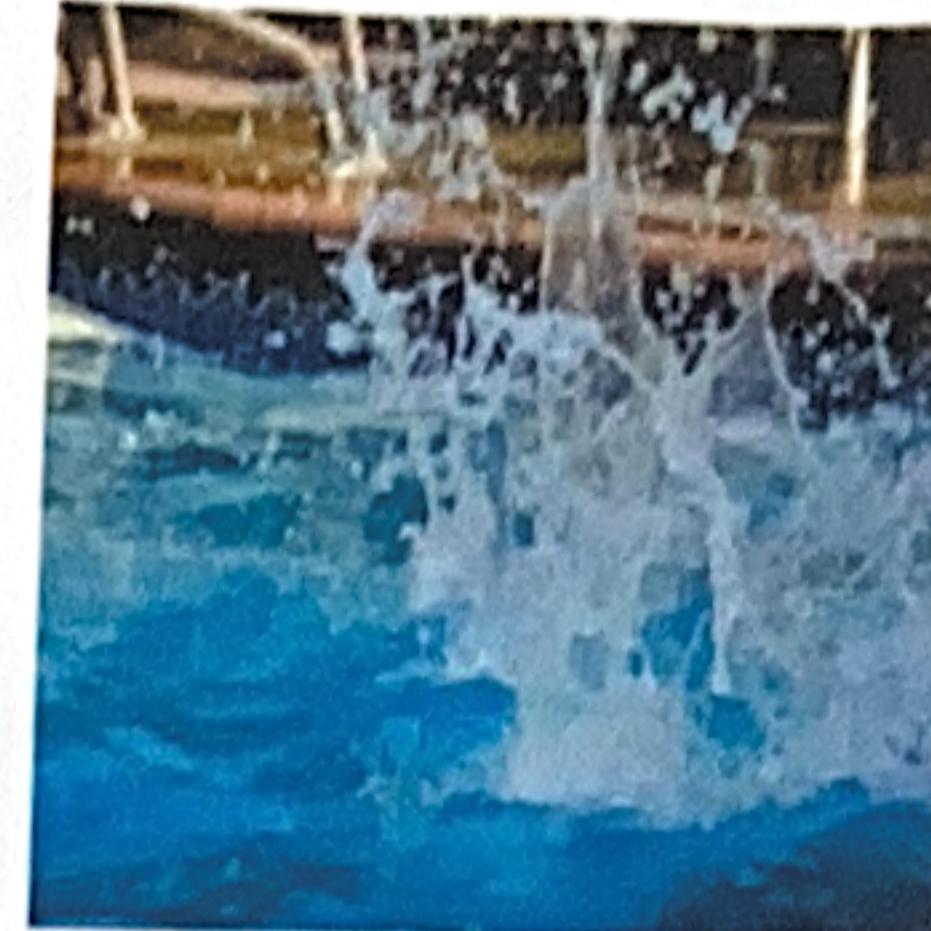


- Fluid Dynamic

Unlike other bodies, fluids exhibit highly volatile behaviors. As a result, it's difficult to come up with a single, efficient way for simulating all fluid effect.



-X: Simulation Approaches

- Lagrangian

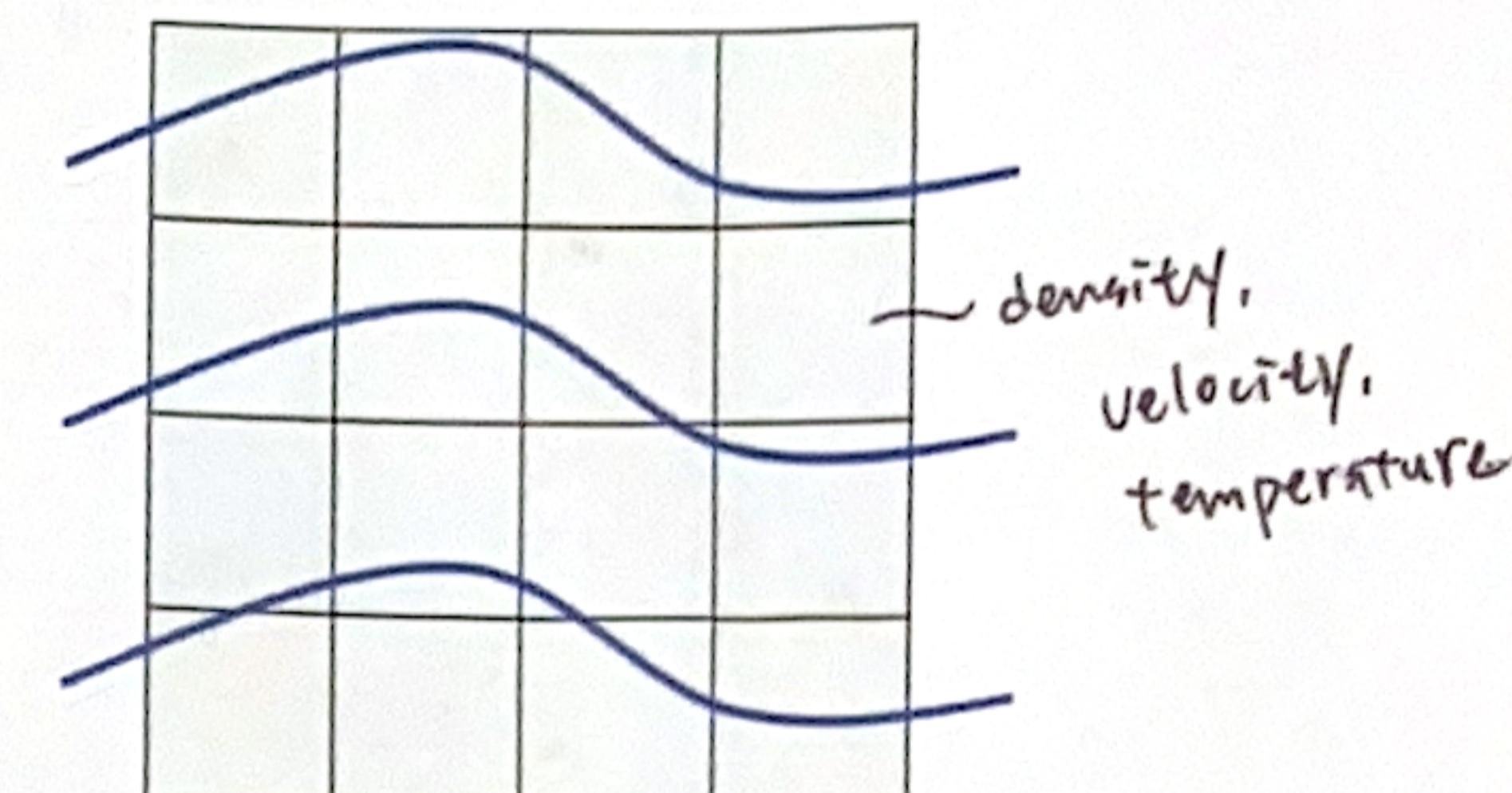
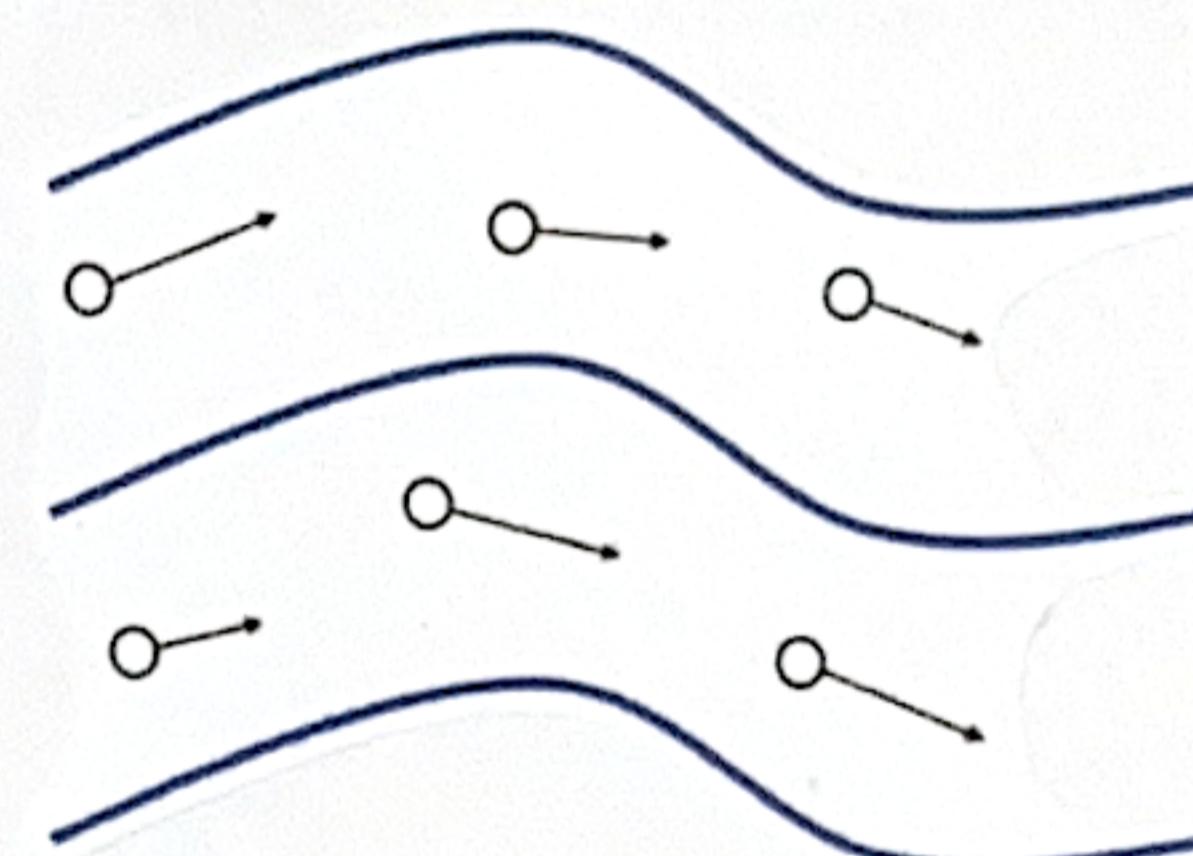
(dynamic particles or mesh)

Node movement carries physical quantities (mass, velocity ...)

- Eulerian

(static grid or mesh)

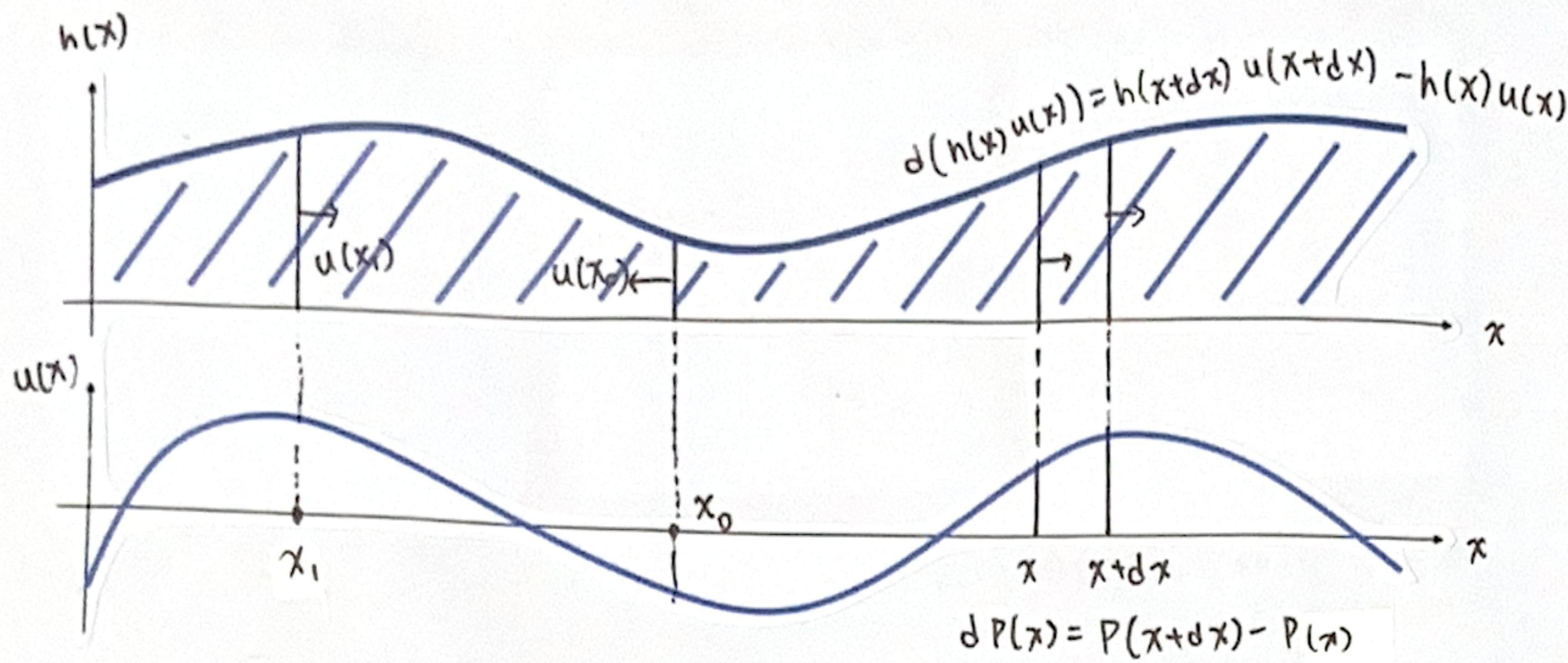
Grid/Mesh doesn't move. Store physical quantities change.



* Height Field Model

- In 2D, a (1.5D) height field is a height function $h(x)$:

$$\frac{dh(x)}{dt} + \frac{d(h(x)u(x))}{dx} = 0$$



- The velocity is also a function $u(x)$: $\frac{du(x)}{dt} = -u(x) \frac{du(x)}{dx} - \frac{1}{\rho} \frac{dP(x)}{dx} + a(x)$

→ ignoring advection and external acceleration, we get a simple form

$$\frac{du(x)}{dt} = -\frac{1}{\rho} \frac{dP(x)}{dx}, \text{ where } \rho \text{ is density and } P(x) \text{ is pressure.}$$

$$\begin{cases} u(x) \downarrow, \text{ when } P(x) > 0 \\ u(x) \uparrow, \text{ when } P(x) < 0 \end{cases}$$

For shallow wave,

$$\left\{ \begin{array}{l} \frac{dh}{dt} + \frac{d(hu)}{dx} = 0 \rightarrow \frac{dh}{dt} + u \frac{dh}{dx} + h \frac{du}{dx} = 0 \xrightarrow{\frac{dt}{dt}} \frac{d^2h}{dt^2} + h \frac{du}{dxdt} = 0 \\ \frac{du}{dt} = -\frac{1}{\rho} \frac{dP}{dx} \xrightarrow{\frac{dx}{dt}} \frac{du}{dt} = -\frac{1}{\rho} \frac{d^2P}{dx^2} \end{array} \right.$$

→ then eliminate u and formulate the shallow wave equation: $\frac{d^2h}{dt^2} = -\frac{1}{\rho} \frac{d^2P}{dx^2}$

$$\rightarrow \frac{h_i^{n+1} - 2h_i^n + h_i^{n-1}}{\Delta t^2} = \frac{h_i^n}{\rho} \frac{P_{i+1}^n - 2P_i^n + P_{i-1}^n}{\Delta x^2} \sim i 表示點, n 表示時間$$

$$\rightarrow h_i^{n+1} = 2h_i^n - h_i^{n-1} + \frac{\Delta t^2 h_i^n}{\Delta x^2 \rho} (P_{i+1}^n - 2P_i^n + P_{i-1}^n)$$

F2

Volume Preservation

The volume of fluid should stay the same when simulation. Suppose that $\sum_i h_i^n = \sum_i h_i^{n+1} = \sum_i h_i^{n-1} = V$.

By summation the difference equation,

$$\begin{aligned} \sum_i h_i^{n+1} &= 2 \sum_i h_i^n - \sum_i h_i^{n-1} + \sum_i \frac{\Delta t^2 h_i^n}{\Delta x^2 \rho} (P_{i+1}^n - 2P_i^n + P_{i-1}^n) \\ &= V + \sum_i \frac{\Delta t^2 h_i^n}{\Delta x^2 \rho} (P_{i+1}^n - 2P_i^n + P_{i-1}^n) \end{aligned}$$

may not be zero

- Solution 1: modify difference equation to

$$h_i^{n+1} = 2h_i^n - h_i^{n-1} + \frac{\Delta t^2}{\Delta x^2 \rho} \left(\left(\frac{h_{i-1}^n + h_i^n}{2} \right) (P_{i-1}^n - P_i^n) + \left(\frac{h_{i+1}^n + h_i^n}{2} \right) (P_{i+1}^n - P_i^n) \right)$$

$$\xrightarrow{\text{sum}} \sum_i h_i^{n+1} = V + \frac{\Delta t^2}{\Delta x^2 \rho} \sum_i \left(\left(\frac{h_{i-1}^n + h_i^n}{2} \right) (P_{i-1}^n - P_i^n) + \left(\frac{h_{i+1}^n + h_i^n}{2} \right) (P_{i+1}^n - P_i^n) \right)$$

This is because water exchanges between h_i^n and h_{i+1}^n .

用 sum 展開後前後項會互相抵消

- Solution 2: assume h_i^n is constant, so

$$h_i^{n+1} = 2h_i^n - h_i^{n-1} + \frac{\Delta t^2 H}{\Delta x^2 \rho} (P_{i+1}^n - 2P_i^n + P_{i-1}^n)$$

$$\xrightarrow{\text{sum}} \sum_i h_i^{n+1} = V + \frac{\Delta t^2 H}{\Delta x^2 \rho} \sum_i ((P_{i-1}^n - P_i^n) + (P_{i+1}^n - P_i^n))$$

must be zero

► Pressure: related to the water height $P_i^n = \rho g h_i^n$

$$\rightarrow h_i^{n+1} = 2h_i^n - h_i^{n-1} + \frac{\Delta t^2 H g}{\Delta x^2} (h_i^n - 2h_i^{n-1} + 2h_{i-1}^n)$$

replace by constant α

► Viscosity: like damping, trying to slow down the wave.

$$\rightarrow h_i^{n+1} = h_i^n + \beta (h_i^n - h_i^{n-1}) + \alpha (h_{i+1}^n - 2h_i^n + 2h_{i-1}^n)$$

momentum

viscosity constant

Shallow wave simulator

for every cell i

$$h_i^{\text{new}} = h_i^{\text{old}} + \beta(h_i - h_i^{\text{old}}) + \alpha(h_{i-1} - h_i) + \alpha(h_{i+1} - h_i)$$

for every cell

$$h_i^{\text{old}} = h_i^{\text{new}}$$

$$h_i = h_i^{\text{new}}$$

Boundary Conditions

• Dirichlet B.C.: boundary height H_{ext} is constant.

It is considered as an open boundary.

• Neumann B.C.: specifies the boundary derivative.

It is considered as a closed boundary

Shallow wave simulator with Neumann B.C.

for every cell i

$$h_i^{\text{new}} = h_i^{\text{old}} + \beta(h_i - h_i^{\text{old}})$$

if h_{i-1} exists, then $h_i^{\text{new}} = h_i^{\text{old}} + \alpha(h_{i-1} - h_i)$

If h_{i+1} exists, then $h_i^{\text{new}} = h_i^{\text{old}} + \alpha(h_{i+1} - h_i)$

for every cell i

$$h_i^{\text{old}} = h_i^{\text{new}}$$

$$h_i = h_i^{\text{new}}$$

Shallow wave simulator with Neumann B.C. in 3D

for every cell i, j

$$h_{i,j}^{\text{new}} = h_{i,j}^{\text{old}} + \beta(h_{i,j} - h_{i,j}^{\text{old}})$$

$$h_{i-1,j}$$

if $h_{i+1,j}$ exists, then $h_{i,j}^{\text{new}} = h_{i,j}^{\text{old}} + \alpha(h_{i+1,j} - h_{i,j})$

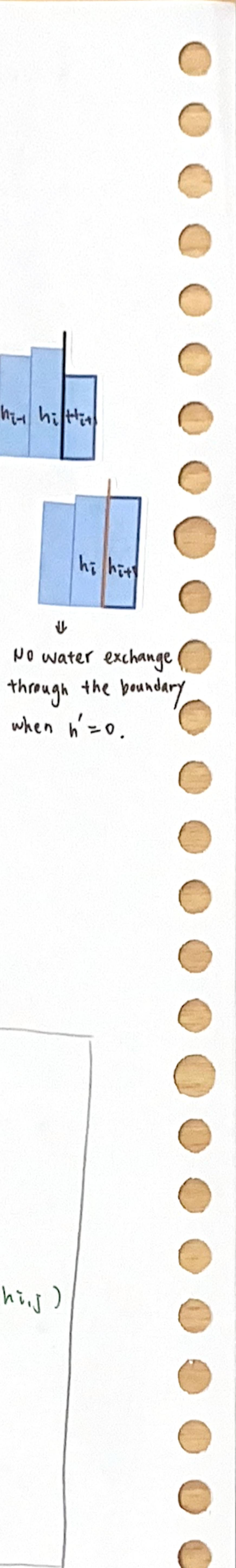
$$h_{i,j-1}$$

$$h_{i,j+1}$$

for every cell i, j

$$h_{i,j}^{\text{old}} = h_{i,j}^{\text{new}}$$

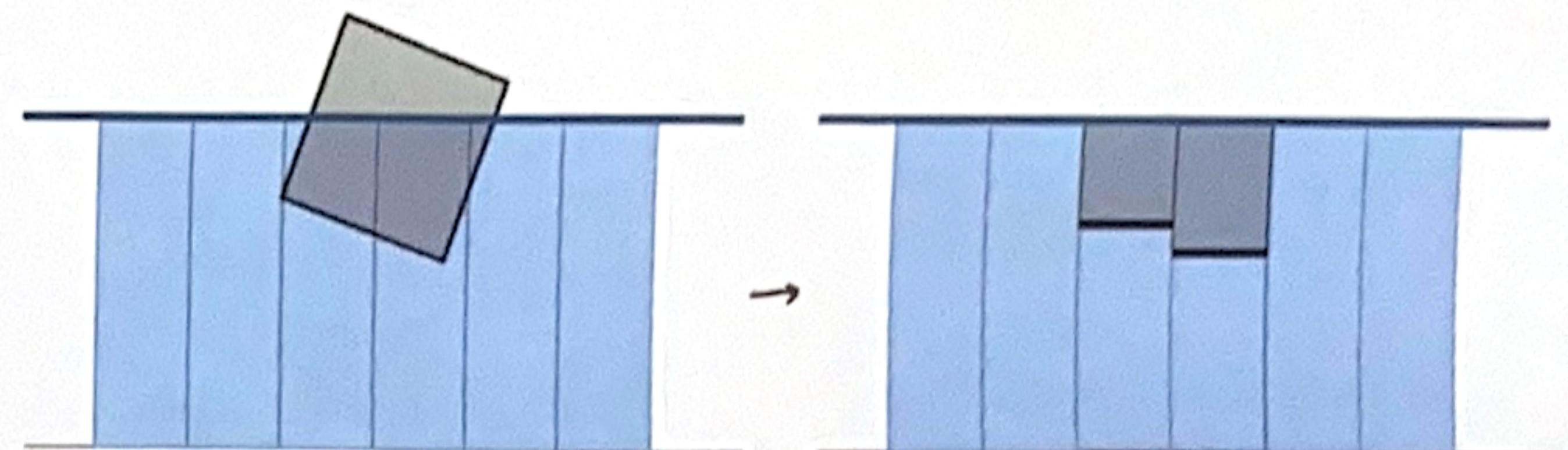
$$h_{i,j} = h_{i,j}^{\text{new}}$$



T3

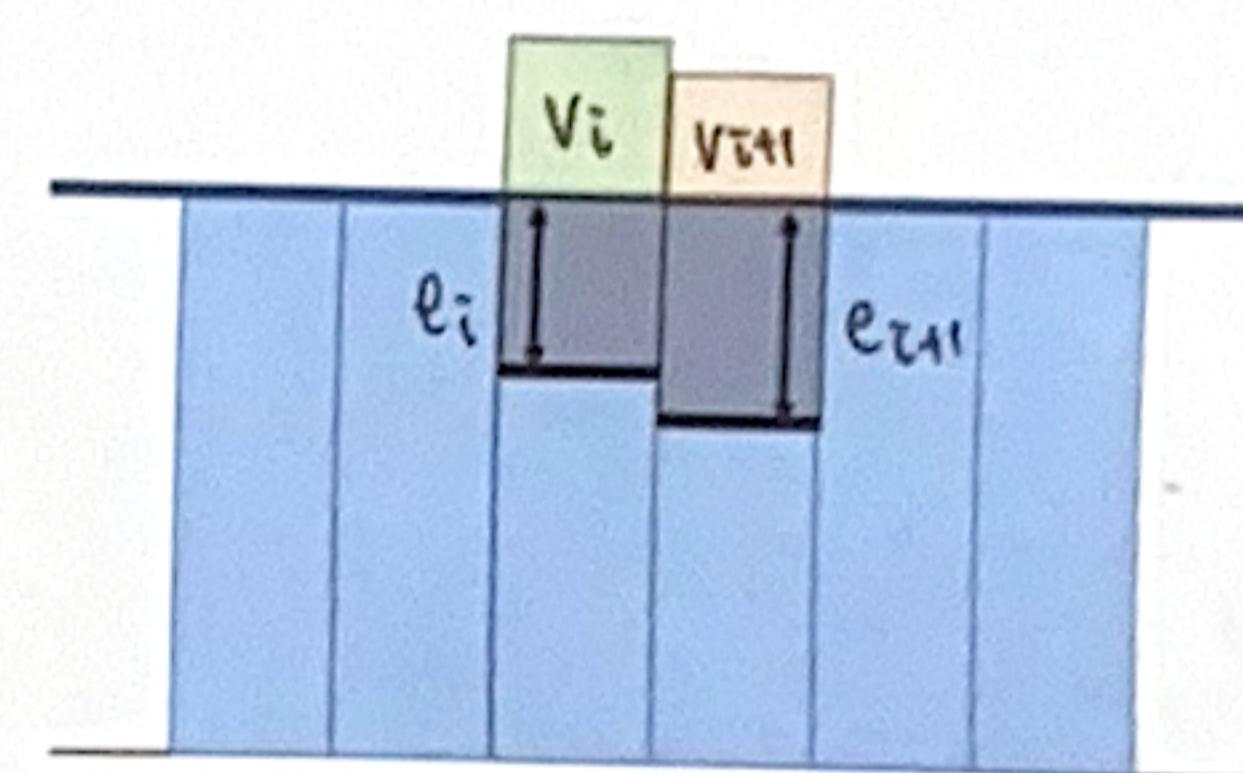
* Two-Way Coupling

- The coupling between a solid and a liquid should be two way, i.e., liquid \rightarrow solid and solid \rightarrow liquid.



How to expel water out of the gray cell regions?

Set up a virtual height v_i , so that $h_i^{\text{real-new}} = h_i - e_i$



$$\rightarrow \begin{cases} h_i - e_i = h_i + \beta(h_i - h_i^{\text{old}}) + \alpha(h_{i+1} + h_{i-1} + h_i - 2v_i - 2h_i) = h_i^{\text{new}} + \alpha(v_{i+1} - 2v_i) \\ \text{modify by } v \end{cases}$$

$$h_{i+1} - e_{i+1} = h_{i+1} + \beta(h_{i+1} - h_{i+1}^{\text{old}}) + \alpha(h_{i+2} + v_i + h_i - 2v_{i+1} - 2h_{i+1}) = h_{i+1}^{\text{new}} + \alpha(v_i - 2v_{i+1})$$

$$\rightarrow \begin{cases} 2v_i - v_{i+1} = \frac{1}{\alpha}(h_i^{\text{new}} - h_i + e_i) = b_i \\ -v_i + 2v_{i+1} = \frac{1}{\alpha}(h_{i+1}^{\text{new}} - h_{i+1} + e_{i+1}) = b_{i+1} \end{cases}$$

$$\rightarrow \begin{bmatrix} 2 & -1 \\ -1 & 2 \end{bmatrix} \begin{bmatrix} v_i \\ v_{i+1} \end{bmatrix} = \begin{bmatrix} b_i \\ b_{i+1} \end{bmatrix}$$

\rightarrow more general in programming

$$\begin{bmatrix} 1 & & & & \\ -1 & 2 & -1 & & \\ & -1 & 2 & -1 & \\ & & -1 & 2 & -1 \\ & & & -1 & 2 \end{bmatrix} \begin{bmatrix} v_{i-1} \\ v_i \\ v_{i+1} \\ v_{i+2} \end{bmatrix} = \begin{bmatrix} 0 \\ b_i \\ b_{i+1} \\ 0 \end{bmatrix}$$

Shallow wave simulator with solid object in 3D

for every cell i, j

$$h_{i,j}^{\text{new}} = h_{i,j} + \beta(h_{i,j} - h_{i,j}^{\text{old}})$$

$h_{i-1,j}$

if $h_{i+1,j}$ exists, then $h_{i,j}^{\text{new}} = h_{i,j}^{\text{new}} + \alpha(h_{i+1,j} - h_{i,j})$

$h_{i,j-1}$

$h_{i,j+1}$

$h_{i,j}$

$h_{i+1,j}$

$h_{i,j-1}$

$h_{i,j+1}$

// Get v

for every cell i, j

if in contact with solid

$$b_{i,j} = \frac{1}{2}(h_{i,j}^{\text{new}} - h_{i,j} + e_{i,j})$$

$\text{tag}_{i,j} = \text{true}$

else $v_{i,j} = 0$

$\text{tag}_{i,j} = \text{false}$

PCG-solver(v, b, tag) \rightarrow 共轭梯度法

for every cell i, j .

$h_{i-1,j}$

if $h_{i+1,j}$ exists, then $h_{i,j}^{\text{new}} = h_{i,j}^{\text{new}} + \alpha(h_{i+1,j} - h_{i,j})$

$h_{i,j-1}$

$h_{i,j+1}$

$h_{i,j}^{\text{old}} = h_{i,j}$

$h_{i,j} = h_{i,j}^{\text{new}}$

relaxation factor, 物体移動
物体太靠近邊界時
浪會很大, because
 $v_{i-1,j}$ $v_{i+1,j}$ $v_{i,j-1}$ $v_{i,j+1}$ this is explicit
method

Rigid Body Update

Estimate the floating force by the actual water expelled in every column by the buoyancy explained by Archimede's principle.

In 2D:

$$f_r = \rho g \Delta x (h_i - h_i^{\text{new}})$$

把物体下移有 Δx ,
 ΔA 造成的浮力加回 F_r

In 3D:

$$f_r = \rho g \Delta A (h_{i,j} - h_{i,j}^{\text{new}})$$

T4

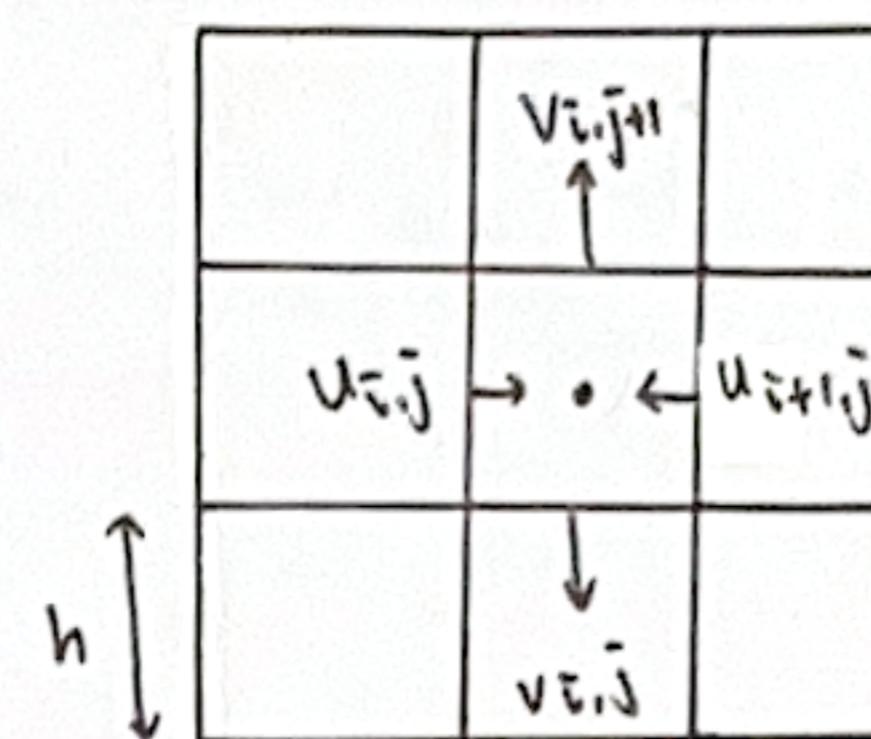
* Incompressible, Viscous Navier-Stokes Equations

Equation formula: { Incompressibility: $\nabla \cdot \vec{u} = 0$

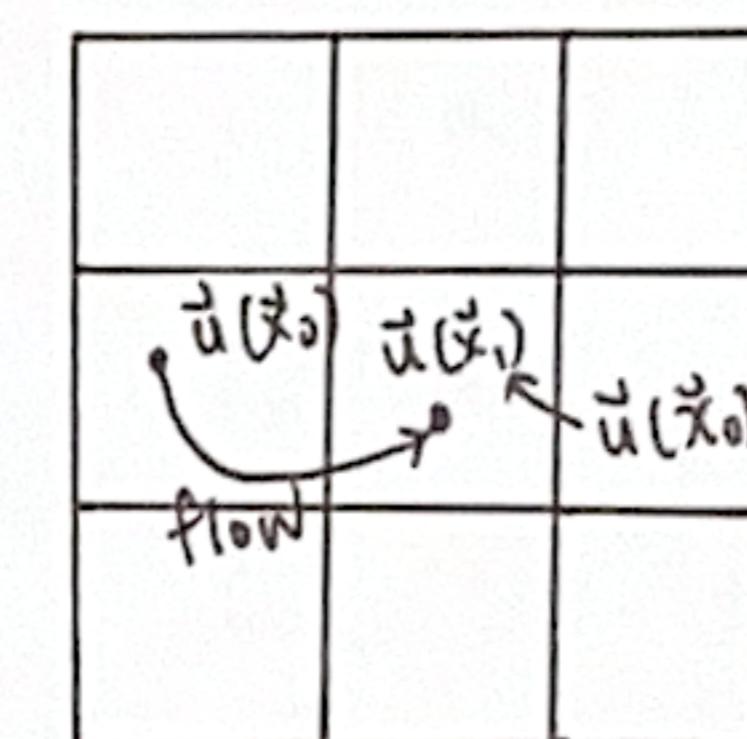
Momentum: $\frac{\partial \vec{u}}{\partial t} = \vec{g} - (\vec{u} \cdot \nabla) \vec{u} + \mu \nabla^2 \vec{u} - \nabla P_{\text{internal}}$
 ↓ Laplacian
 ↓ advection diffusion
 external acceleration
 (ex. gravity)

Using method of characteristics to solve a long partial differential equation in steps.

• Step 1: external acceleration

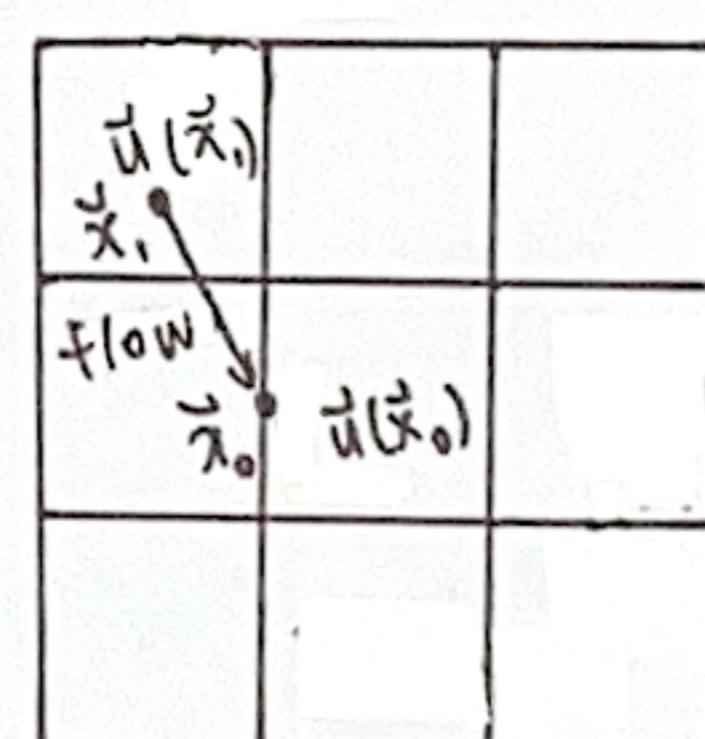


• Step 2: advection



To solve this problem, we come to realize that advection means to carry physical quantities by velocity. The solution is to trace a virtual particle backward over time. The Semi-Lagrangian Method:

a. One step:



b. Three steps:

compute $u(x_0)$

(take x-direction
for example)

$$x_1 = x_0 - \Delta t u(x_0)$$

compute $u(x_1)$

$$u_{i,j}^{\text{new}} = u(x_1)$$

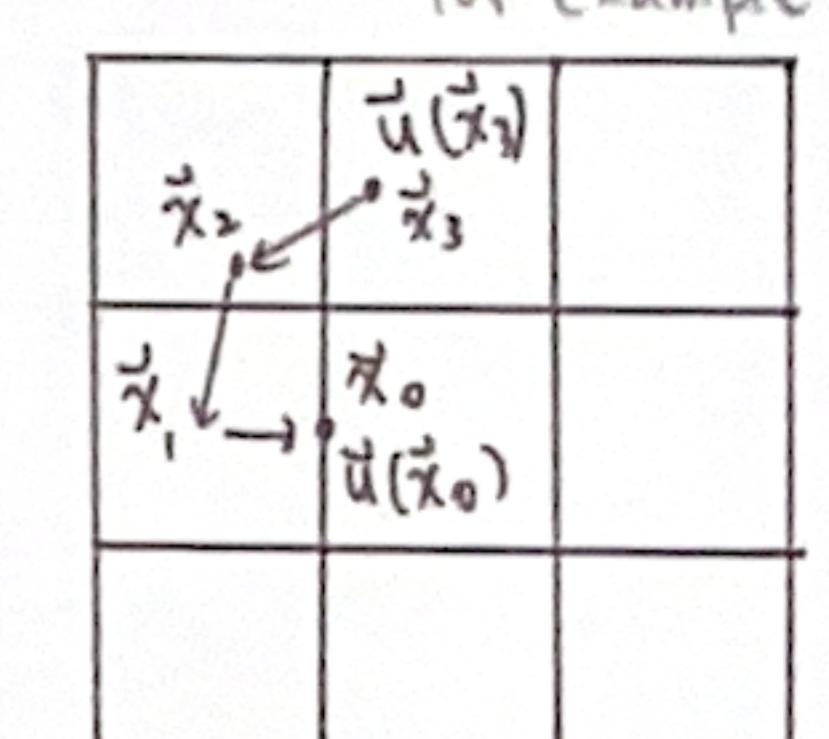
Define $y_0 = (i, j - 0.5)$

compute $v(y_0)$

$$y_1 = y_0 - \Delta t v(y_0)$$

compute $v(y_1)$

$$v_{i,j}^{\text{new}} = v(y_1)$$



" \rightarrow " means flow

Define $x_0 = (i - 0.5, j)$
compute $u(x_0)$

$$x_1 = x_0 - \frac{1}{3} \Delta t u(x_0)$$

compute $u(x_1)$

$$x_2 = x_1 - \frac{1}{3} \Delta t u(x_1)$$

compute $u(x_2)$

$$x_3 = x_2 - \frac{1}{3} \Delta t u(x_2)$$

compute $u(x_3)$

$$u_{i,j}^{\text{new}} = u(x_3)$$

• Step 3: diffusion

Update \vec{u} by solving $\frac{\partial \vec{u}}{\partial t} = \mu \Delta \vec{u}$

a. One step:

$$u_{i,j}^{\text{new}} = u_{i,j} + \mu \Delta t \frac{u_{i-1,j} + u_{i+1,j} + u_{i,j-1} + u_{i,j+1}}{h^2}$$

$$v_{i,j}^{\text{new}} = v_{i,j} + \mu \Delta t \frac{v_{i-1,j} + v_{i+1,j} + v_{i,j-1} + v_{i,j+1}}{h^2}$$

If $\mu \Delta t$ is large, the system would be unstable, so we could also use smaller sub-steps.

→ b. Two steps: (take x -direction for example)

$$u_{i,j}^{\text{temp}} = u_{i,j} + \mu \frac{\Delta t}{2} \frac{u_{i-1,j} + u_{i+1,j} + u_{i,j-1} + u_{i,j+1}}{h^2}$$

$$u_{i,j}^{\text{new}} = u_{i,j}^{\text{temp}} + \mu \frac{\Delta t}{2} \frac{u_{i-1,j}^{\text{temp}} + u_{i+1,j}^{\text{temp}} + u_{i,j-1}^{\text{temp}} + u_{i,j+1}^{\text{temp}}}{h^2}$$

• Step 4: Pressure projection

p on grid!

$p_{i,j}$	$p_{i,j}$	
	$p_{i,j}$	

Update \vec{u} by solving $\frac{\partial \vec{u}}{\partial t} = -\nabla p$

Staggered grid makes this easy:

$$\begin{cases} u_{i,j}^{\text{new}} = u_{i,j} - \frac{1}{h} (p_{i,j} - p_{i-1,j}) \\ v_{i,j}^{\text{new}} = v_{i,j} - \frac{1}{h} (p_{i,j} - p_{i,j-1}) \end{cases}$$

What is p ? It is the pressure caused by incompressibility, which means that after updating from this step, the velocity field should achieve

$$\nabla \cdot \vec{u}^{\text{new}} = 0$$

which means $u_{i+1,j}^{\text{new}} + v_{i,j+1}^{\text{new}} - u_{i,j}^{\text{new}} - v_{i,j}^{\text{new}} = 0$ (4 walls' velocity on a grid)

$$\rightarrow u_{i+1,j} - \frac{1}{h} (p_{i+1,j} - p_{i,j}) + v_{i,j+1} - \frac{1}{h} (p_{i,j+1} - p_{i,j})$$

$$- u_{i,j} + \frac{1}{h} (p_{i,j} - p_{i-1,j}) - v_{i,j} + \frac{1}{h} (p_{i,j} - p_{i,j-1}) = 0$$

$$\rightarrow 4p_{i,j} - p_{i+1,j} - p_{i,j+1} - p_{i-1,j} - p_{i,j-1} = h(u_{i,j} + v_{i,j} - u_{i+1,j} - v_{i,j+1})$$

With Boundary Conditions

$$\begin{cases} \text{Dirichlet (open)} & p_{i-1,j} = P \\ \text{Neumann (close)} & p_{i-1,j} = p_{i,j} \end{cases}$$

→ Once we solve p , update \vec{u} and done.

F5

• Air and Smoke

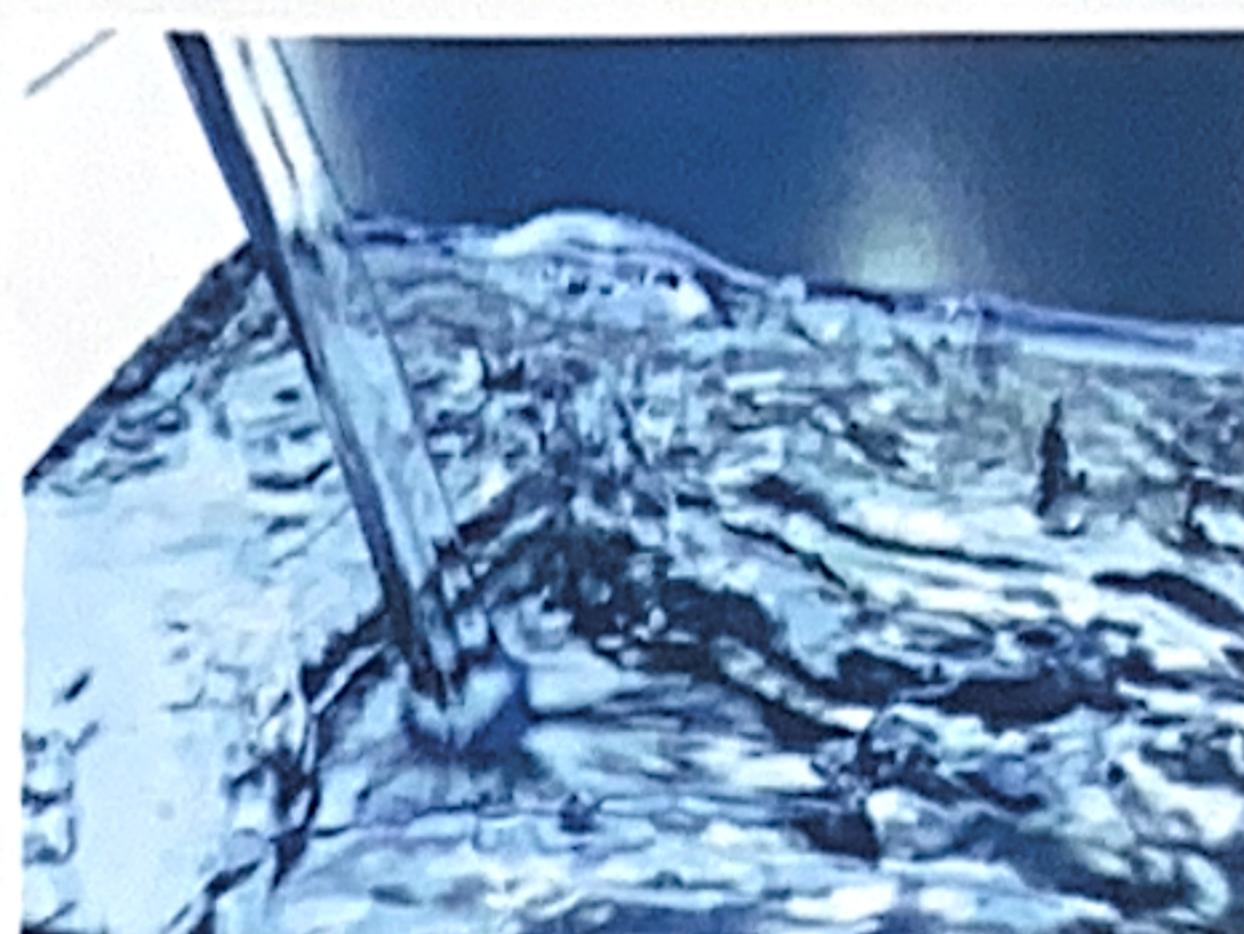


• Air simulation is done in two steps.

1. update the flow (the velocity field) \vec{u} .
2. use semi-Lagrangian advection all of the other physical quantities, i.e., density, temperature.....

- Typically we use Dirichlet boundaries for an open space; Neumann boundaries for a container.
- It can be used to simulate underwater as well.

• Water Simulation



- Two presentations
 - Volume-of-fluid
 - Signed distance function defined over the grid

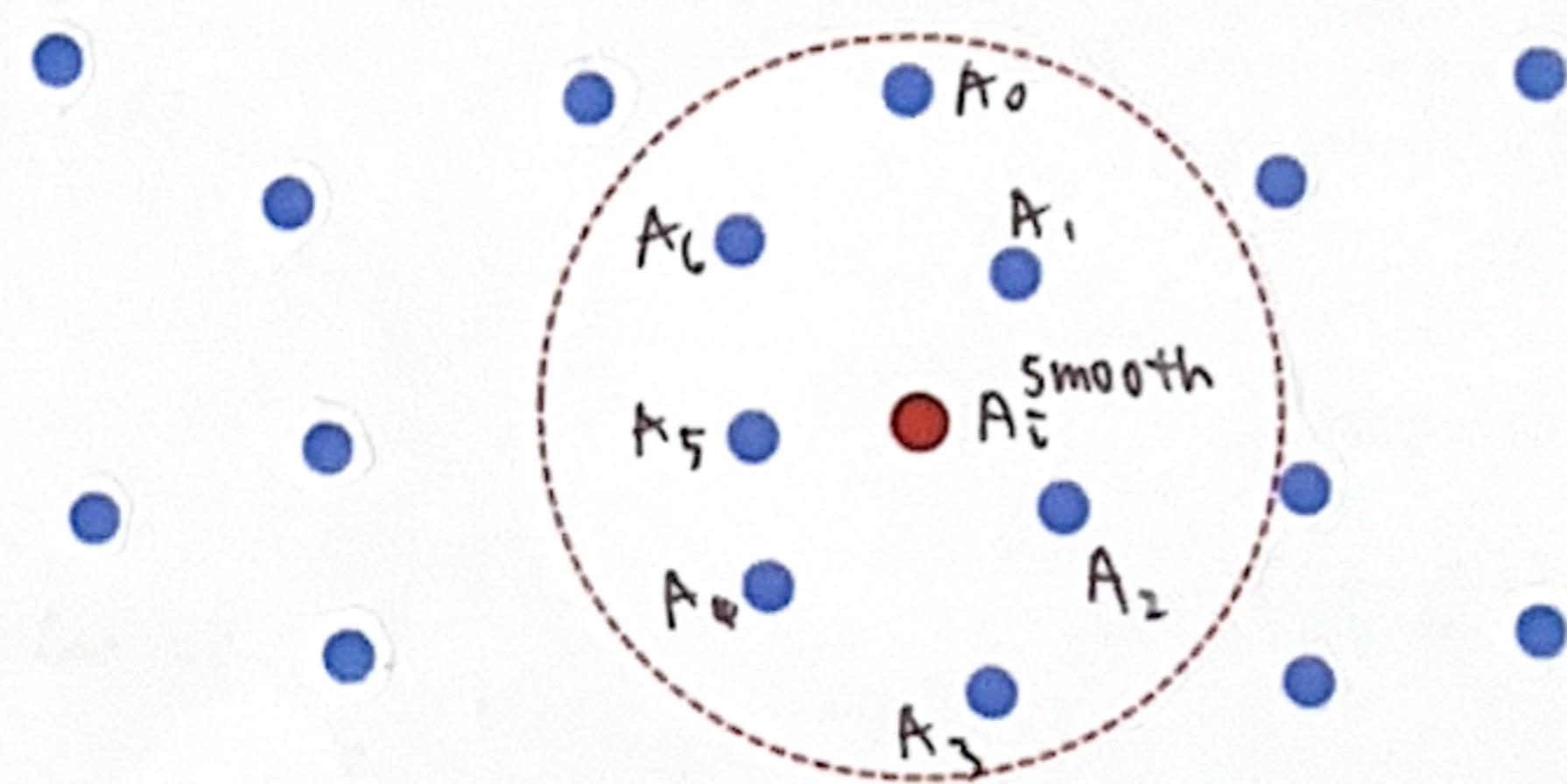
- When implementing advection, both Semi-Lagrangian and level set method cause volume loss. → need corrections.

• Smoothed Particle Hydrodynamics (SPH)

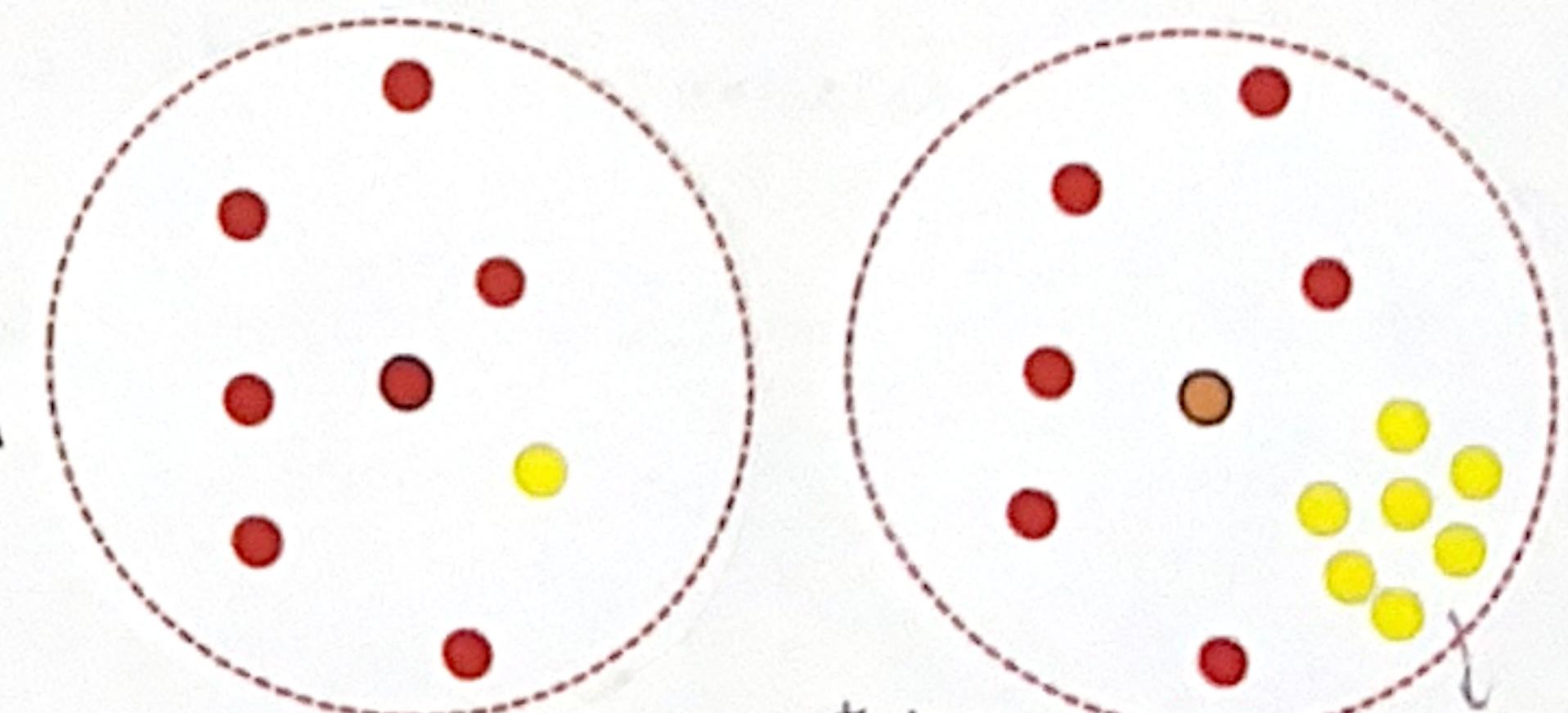
* SPH Model

Consider a Lagrangian particle system: each water molecule is a particle with physical quantities attached, such as position \vec{x}_i , velocity \vec{v}_i , and mass m_i generally write as A_i .

• Simple model: $A_i^{\text{smooth}} = \frac{1}{n} \sum_j A_j$, for $\|\vec{x}_i - \vec{x}_j\| < R$



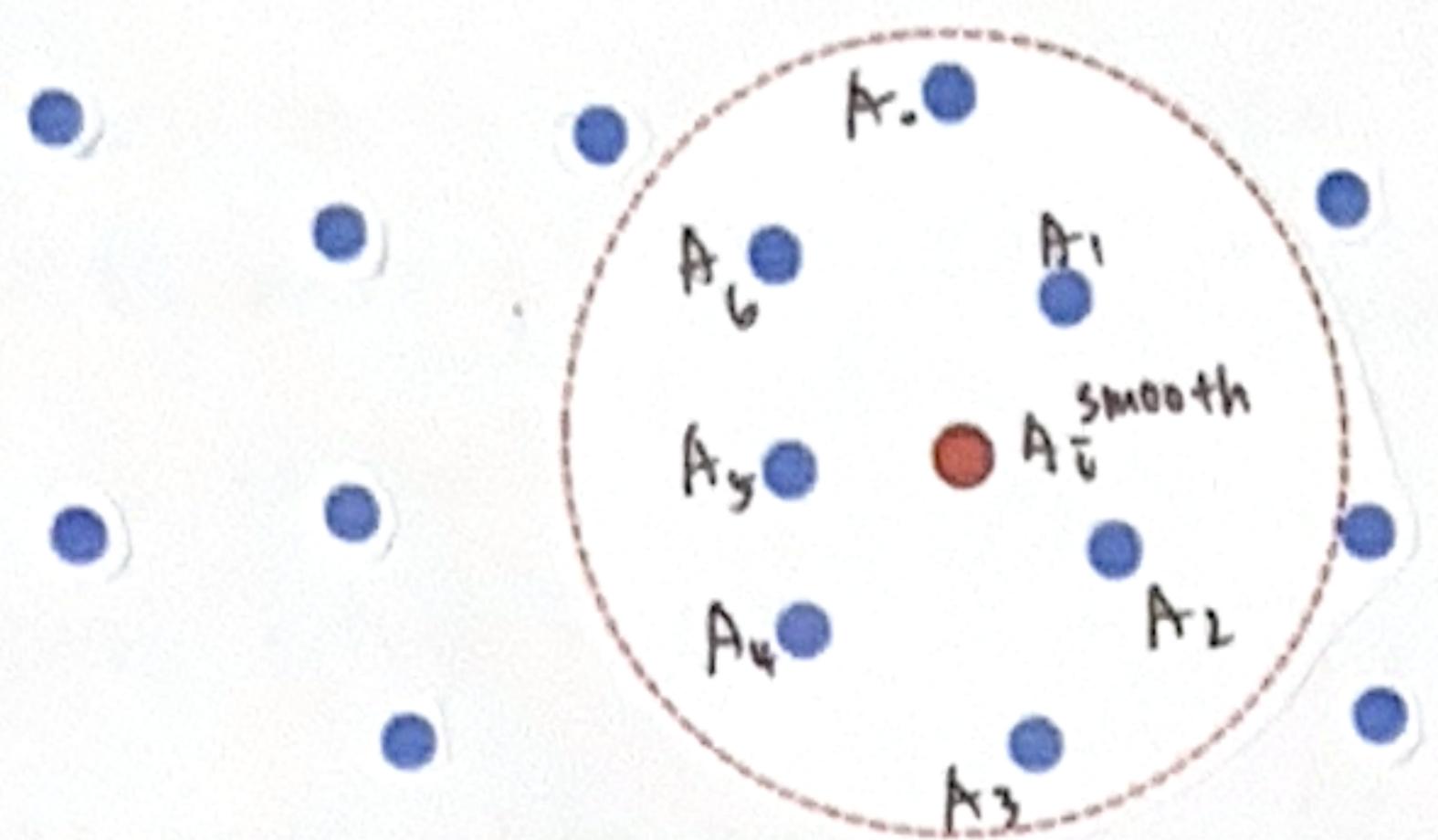
- Problem: only considered the quantity of particles. Didn't consider the distribution of each particle.



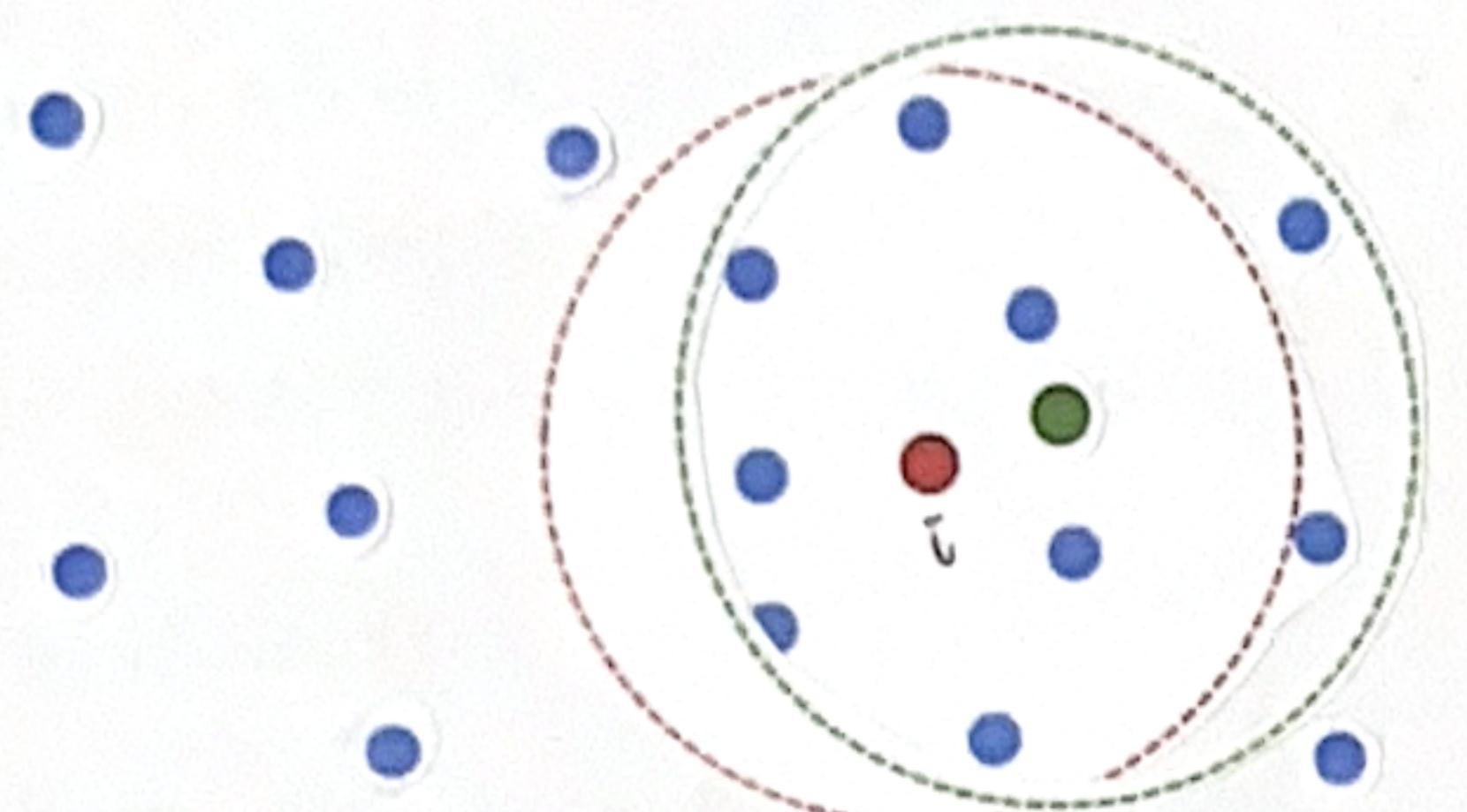
黄色空间小但数量多！

- Better Model: assume each particle has a volume V_j

$$\rightarrow A_i^{\text{smooth}} = \frac{1}{n} \sum_j V_j A_j, \text{ for } \|x_i - x_j\| < R$$



- Problem: not smooth when moving the selection area! ($\rightarrow a)$



Final Model: $A_i^{\text{smooth}} = \sum_j V_j A_j W_{ij}$ for $\|x_i - x_j\| < R$

where W_{ij} is the smoothing kernel. $\begin{cases} \|x_i - x_j\| \text{ large, } W_{ij} \text{ small} \\ \|x_i - x_j\| \text{ small, } W_{ij} \text{ large} \end{cases}$

Volume of particle i : $V_i = \frac{m_i}{\rho_i} \rightarrow V_i = \frac{m_i}{\rho_i^{\text{smooth}}} = \frac{m_i}{\sum_j m_j W_{ij}}$

(By the final model, $\rho_i^{\text{smooth}} = \sum_j V_j \rho_j W_{ij} = \sum_j m_j W_{ij}$)

Therefore, the actual solution is

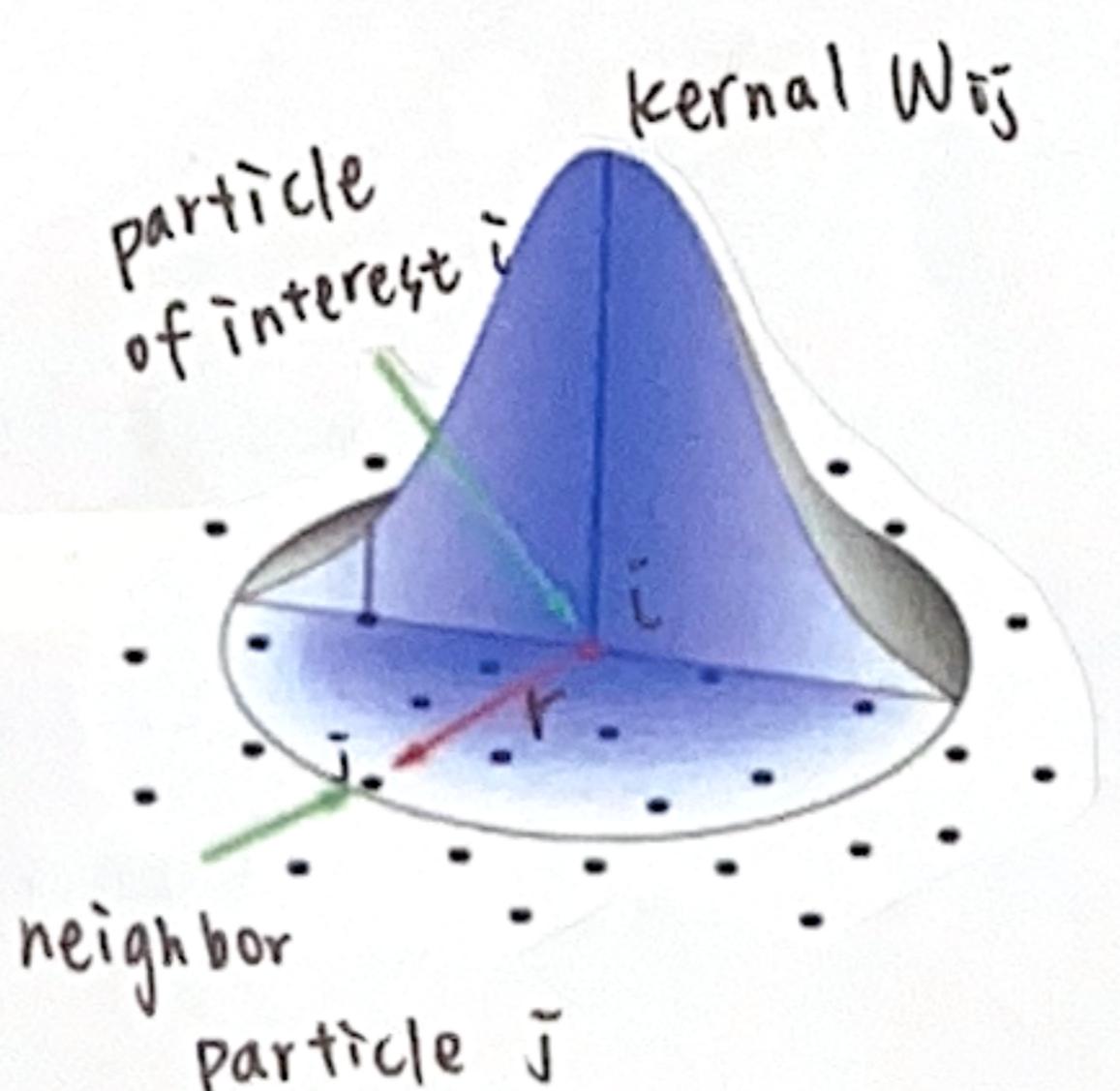
$$A_i^{\text{smooth}} = \sum_j \frac{m_j}{\sum_k m_k W_{ik}} A_j W_{ij}$$

- The derivative is easy to compute:

Gradient: $\nabla A_i^{\text{smooth}} = \sum_j V_j A_j \nabla W_{ij}$

Laplacian: $\nabla^2 A_i^{\text{smooth}} = \sum_j V_j A_j \nabla^2 W_{ij}$

- There are a lot of kernel models to calculate W_{ij} . For example →



F_b

* SPH-Based Fluids

Modeled fluid dynamics by applying three forces on particle i ,

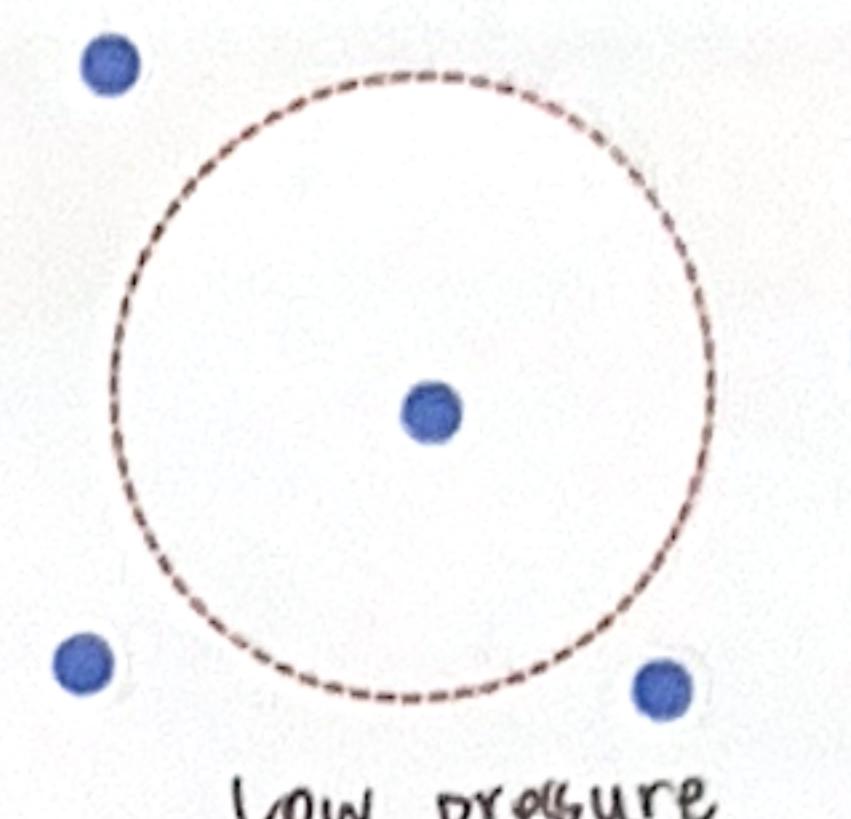
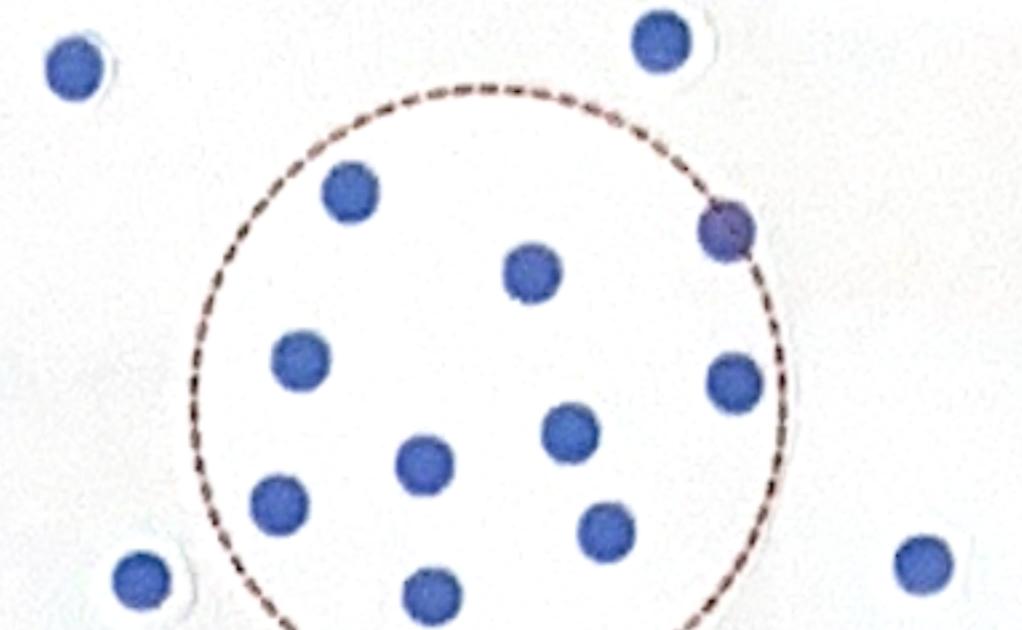
1. Gravity Force: $\vec{F}_i^{\text{gravity}} = m_i g$

2. Pressure Force: related to the density

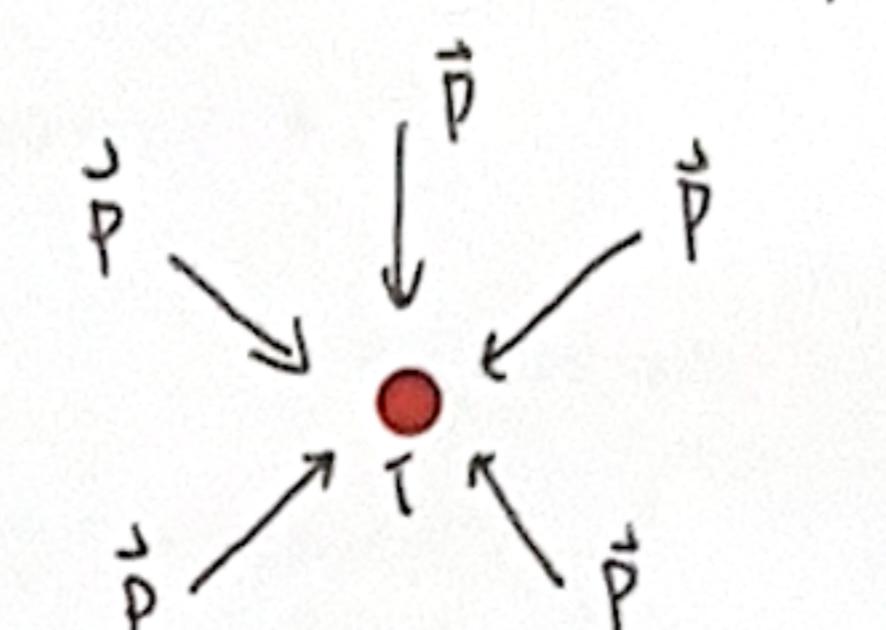
a. compute density of particle i : $\rho_i = \sum_j m_j W_{ij}$

- b. convert it into pressure (by some empirical function):

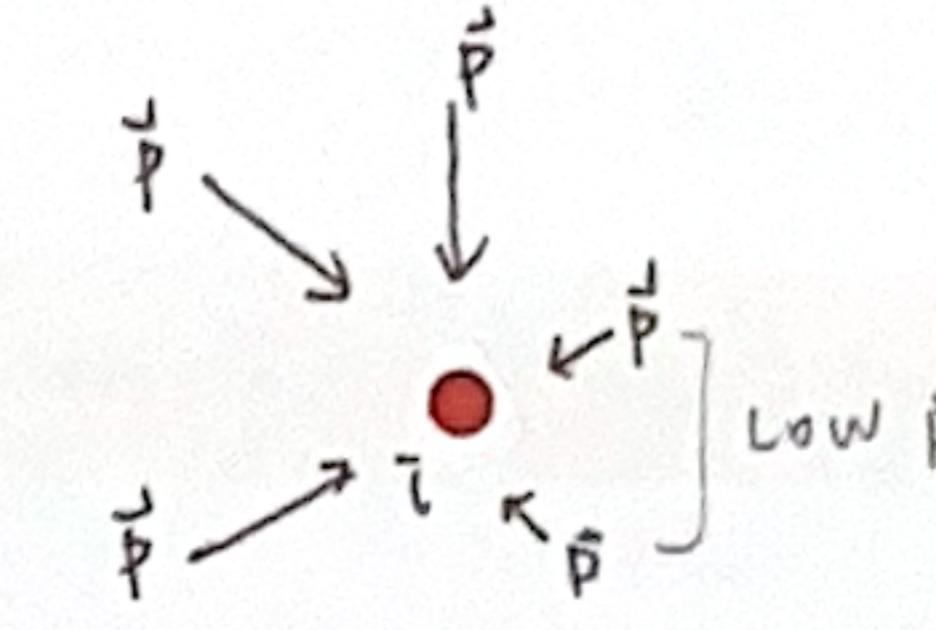
$$\rho_i = k \left(\left(\frac{p_i}{p_{\text{constant}}} \right)^n - 1 \right)$$



- c. Pressure force depends on the difference of pressure:



→ NO pressure force!

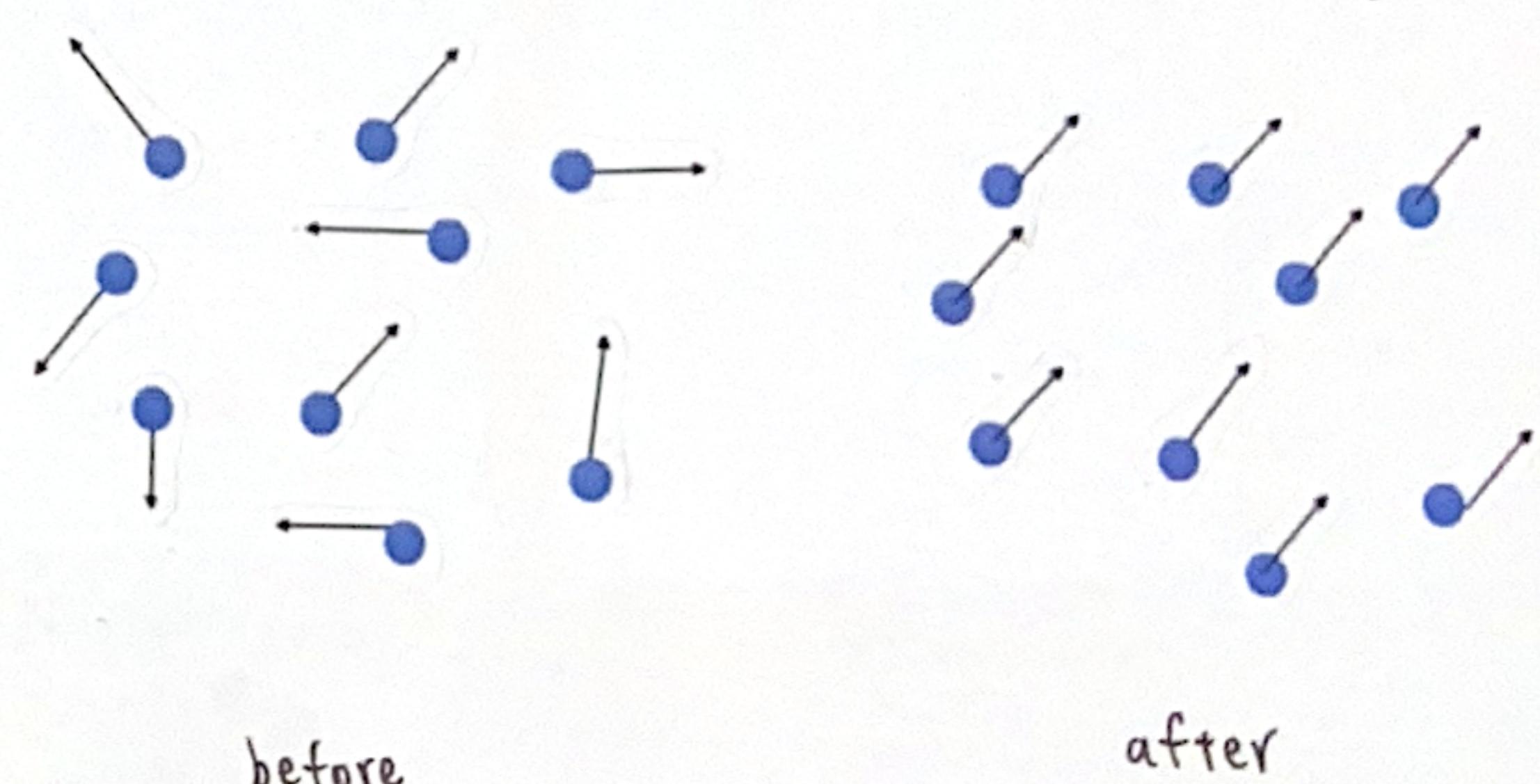


→ Pressure force

Therefore, mathematically, the difference of pressure is gradient of pressure: $\vec{F}_i^{\text{pressure}} = -V_i \nabla_i p^{\text{smooth}} = -V_i \sum_j V_j p_j \nabla_i W_{ij}$

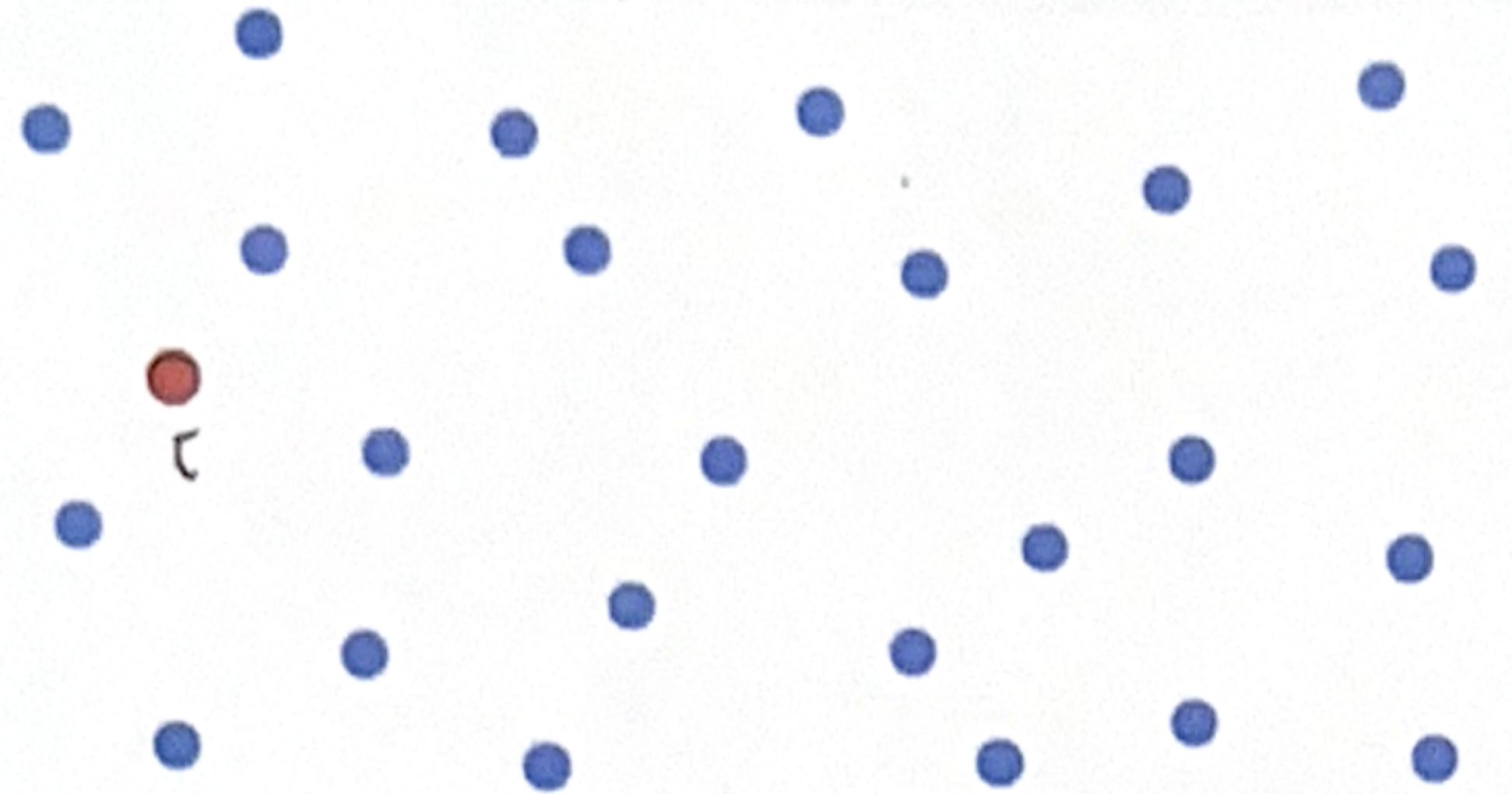
3. Viscosity Force: in order to minimize the difference between the particle velocity and the velocities of its neighbors.

$$\vec{F}_i^{\text{viscosity}} = -V_i m_i \nabla_i^2 v^{\text{smooth}} = -V_i m_i \sum_j V_j \vec{v}_j \nabla_i^2 W_{ij}$$

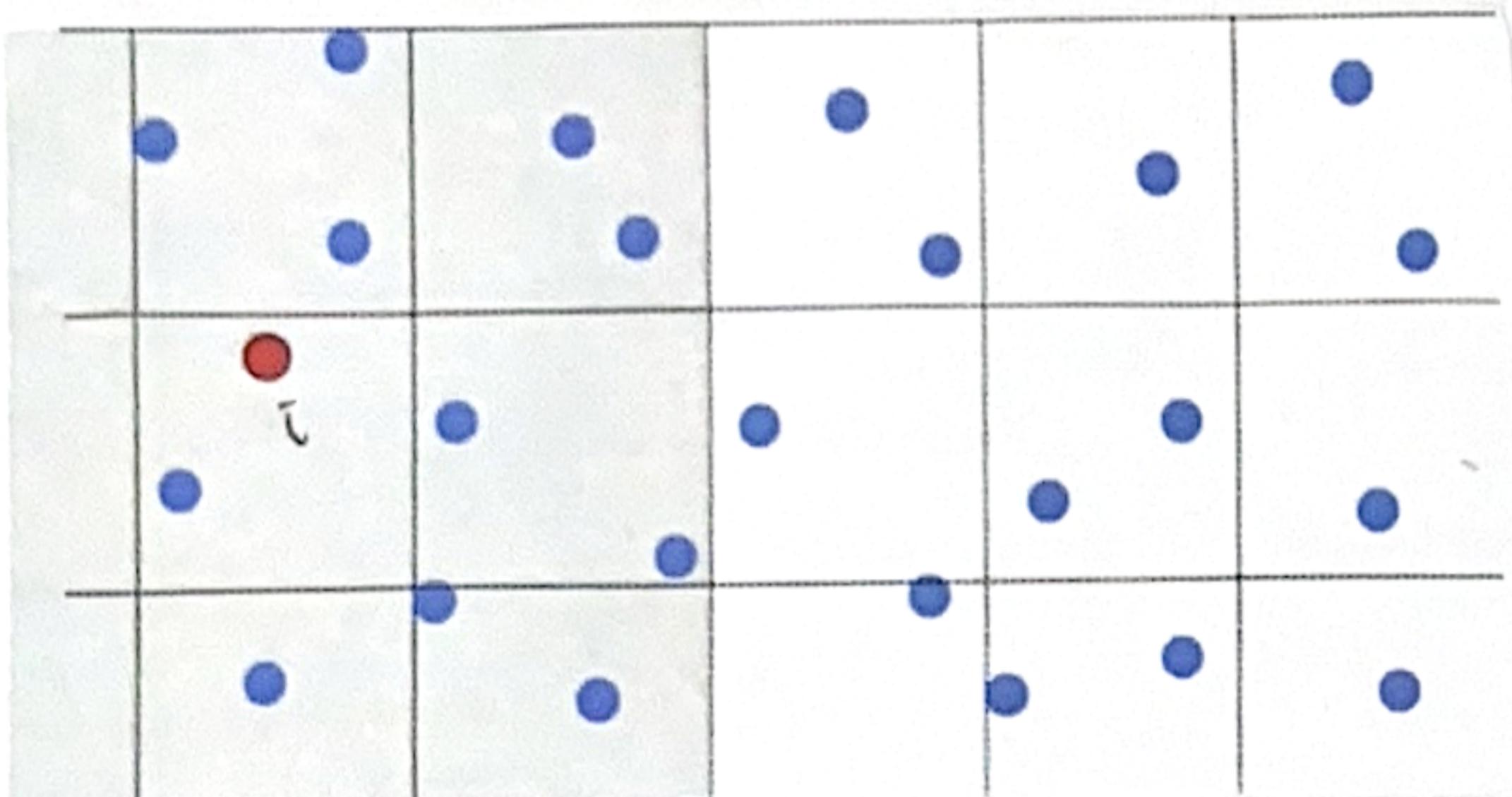


► Problem: exhaustive neighborhood search

To search over every particle pair, the time complexity is $O(N^2)$.

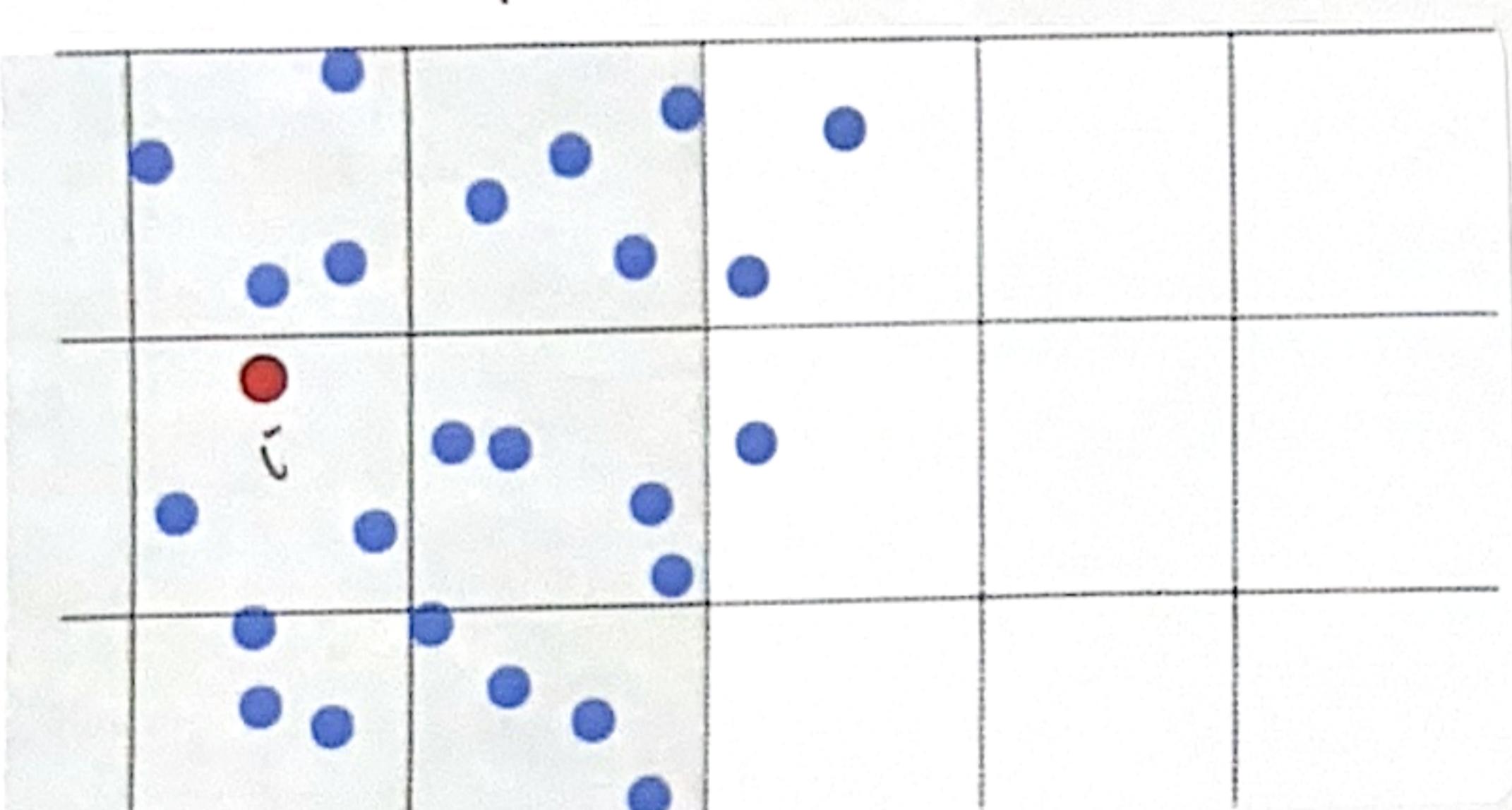


Solution: Spatial Partition



- Separate the space into cell
- Each cell stores the particles in it
- To find the neighborhood of i , just look at the surrounding.

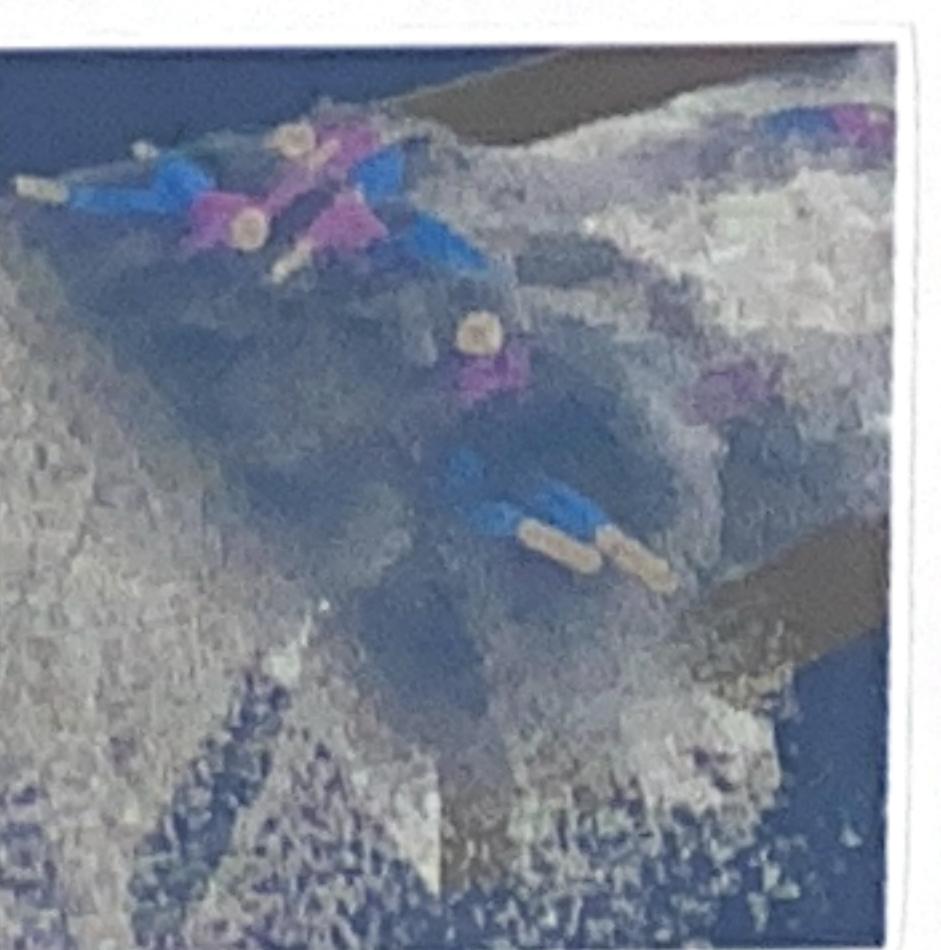
If particles are not uniformly distributed:



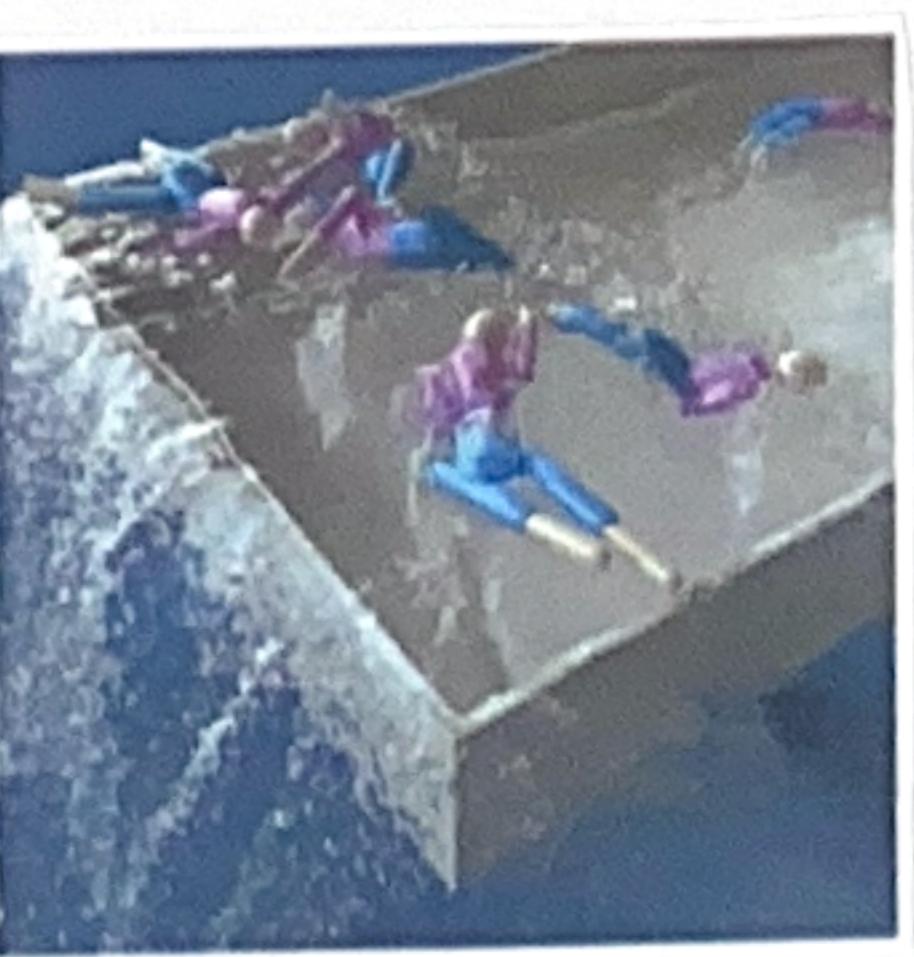
- Octree
- Binary Spatial Partitioning tree

► Fluid display

Reconstruct the water surface from particles.



representation



typical visualization