

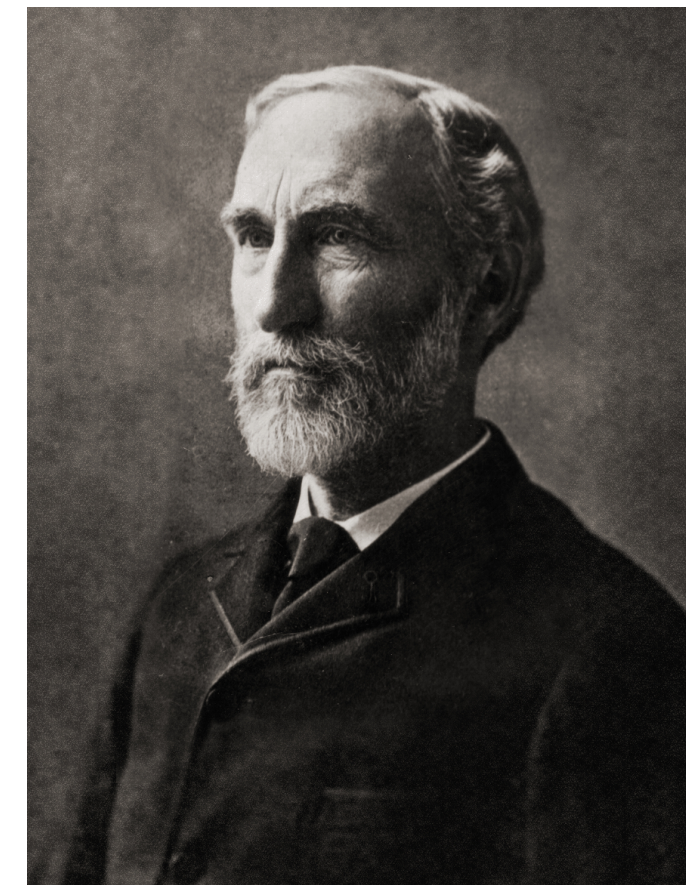
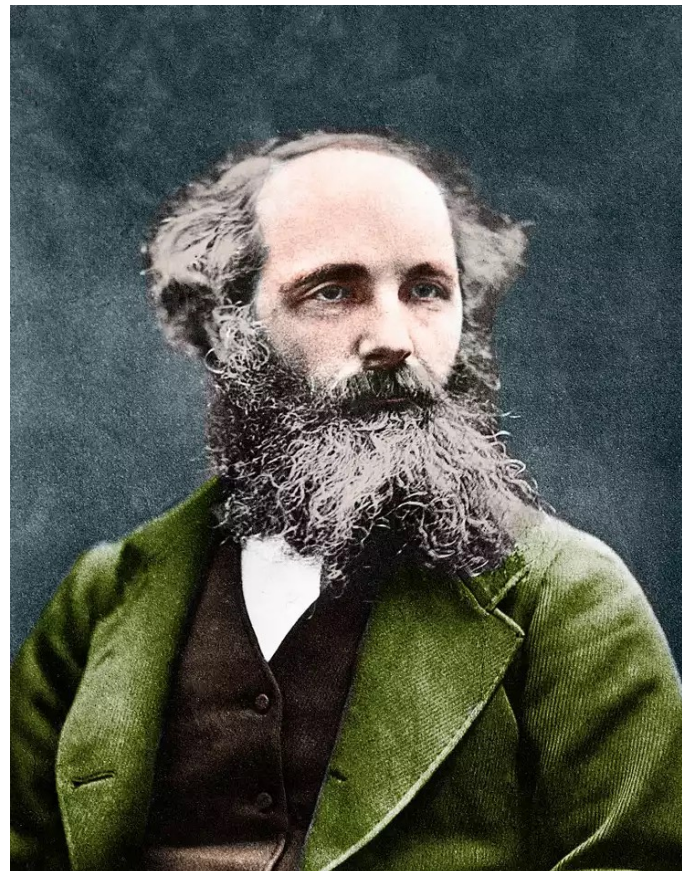
# Random point configurations in statistical physics

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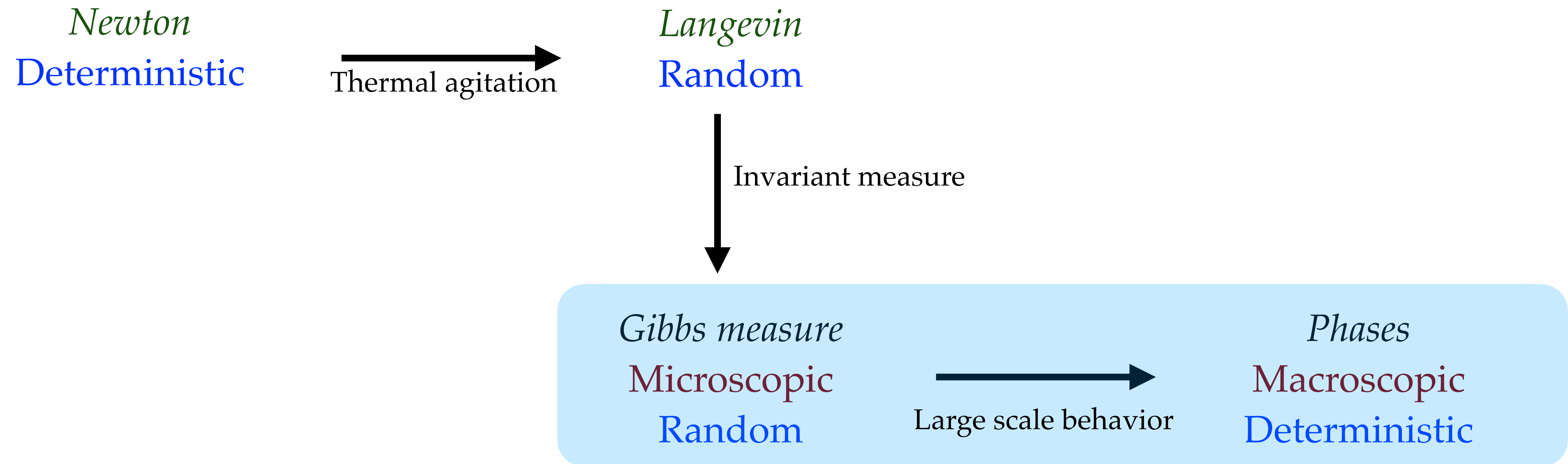
**Point configurations: from statistical physics to potential theory (CIRM, May 2026)**

# Statistical physics

- Statistical physics = physics of **many, many elements in interaction**  
e.g.  $H_2O$  in a pot, molecules in the air of the room...
- Very large number  $N$  of elements (**1 mole =  $10^{23}$** )... almost « infinite » !
- Developed in late 19th / early 20th century (**Maxwell, Boltzmann, Gibbs**)
- *Thermodynamics* was phenomenological, *statistical mechanics* is **reductionist**



# A reductionist approach?



Random distributions (*ensembles*)

# Ingredients of a model

1. *State space*  $S$  (all possible states of the system)  
e.g.  $S = (\mathbb{R}^d)^N$  for  $N$  particles in  $\mathbb{R}^d$
2. Reference *measure*  $\mu$   
e.g.  $\mu =$  Lebesgue, or Riemannian volume on manifold
3. *Energy functional*  $H : S \rightarrow (-\infty, +\infty]$   
e.g.  $H =$  sum of all interactions between particles
4. *Inverse temperature* parameter  $\beta > 0$   
sets the *strength* of interactions

# The basic recipe

The state  $\mathbf{X} \in S$  is *random* and distributed as:

$$d\mathbb{P}(\mathbf{X}) = \frac{1}{Z} e^{-\beta H(\mathbf{X})} d\mu(\mathbf{X})$$

Gibbs measure

Partition function

Boltzmann's factor

The diagram shows the equation  $d\mathbb{P}(\mathbf{X}) = \frac{1}{Z} e^{-\beta H(\mathbf{X})} d\mu(\mathbf{X})$ . Each term is enclosed in a colored oval:  $d\mathbb{P}(\mathbf{X})$  in a reddish-brown oval,  $Z$  in a green oval,  $e^{-\beta H(\mathbf{X})}$  in a blue oval, and  $d\mu(\mathbf{X})$  in a yellow oval. Three arrows point from text labels to these ovals: 'Gibbs measure' points to  $d\mathbb{P}(\mathbf{X})$ , 'Partition function' points to  $Z$ , and 'Boltzmann's factor' points to  $e^{-\beta H(\mathbf{X})}$ .

$$Z = \int_S e^{-\beta H(\mathbf{X})} d\mu(\mathbf{X})$$

# Microstates, macrostates

- $\mathbf{X} \in S$  a *microstate* = a description of the system at **microscopic** scale  
e.g. **all** the positions and **all** the velocities of **all** the molecules in the room  
*precise but not very useful*
- *Observable*  $\mathcal{O} : S \rightarrow Y$  = function of the state, preferably of *macroscopic* nature  
e.g. “the **average** kinetic energy of the molecules in **this corner** of the room”
- *Macrostate* = “all the microstates  $\mathbf{X}$  such that  $\mathcal{O}(\mathbf{X}) =$  a certain value”.
- Goal: understand typical values of chosen  $\mathcal{O}$ , namely *typical macrostates*

*Typical ?? Under the Gibbs measure!*

Gibbs measure = distribution of *microstates*

$$d\mathbb{P}(\mathbf{X}) = \frac{1}{Z} e^{-\beta H(\mathbf{X})} d\mu(\mathbf{X})$$

Law of *observable*

$$\mathbb{P}(\mathcal{O}(\mathbf{X}) = \lambda) = \frac{1}{Z} \int_S e^{-\beta H(\mathbf{X})} \mathbf{1}_{\mathcal{O}(\mathbf{X}) = \lambda} d\mu(\mathbf{X})$$

*Energy*: from " $\mathcal{O}(\mathbf{X}) = \lambda$ ", can I say that  $H(\mathbf{X}) \approx \tilde{H}_\lambda$ ?

$$\mathbb{P}(\mathcal{O}(\mathbf{X}) = \lambda) \approx \frac{1}{Z} e^{-\beta \tilde{H}_\lambda} \int_S \mathbf{1}_{\mathcal{O}(\mathbf{X}) = \lambda} d\mu(\mathbf{X})$$

Can I compute the *volume*

$$\int_S \mathbf{1}_{\mathcal{O}(\mathbf{X}) = \lambda} d\mu(\mathbf{X}) = \mu(\mathcal{O}(\mathbf{X}) = \lambda) ?$$

$$\mathbb{P}(\mathcal{O}(\mathbf{X}) = \lambda) \approx \frac{1}{Z} e^{-\beta \tilde{H}_\lambda} \times \mu(\mathcal{O}(\mathbf{X}) = \lambda)$$

$$\mathbb{P}(\mathcal{O}(\mathbf{X}) = \lambda) = \frac{1}{Z} e^{-\beta \tilde{H}_\lambda + \log \mu(\mathcal{O}(\mathbf{X}) = \lambda)}$$

# Free energy

$$\mathbb{P}(\mathcal{O}(\mathbf{X}) = \lambda) \approx \frac{1}{Z} e^{-\beta \tilde{H}_\lambda + \log \mu(\mathcal{O}(\mathbf{X}) = \lambda)}$$

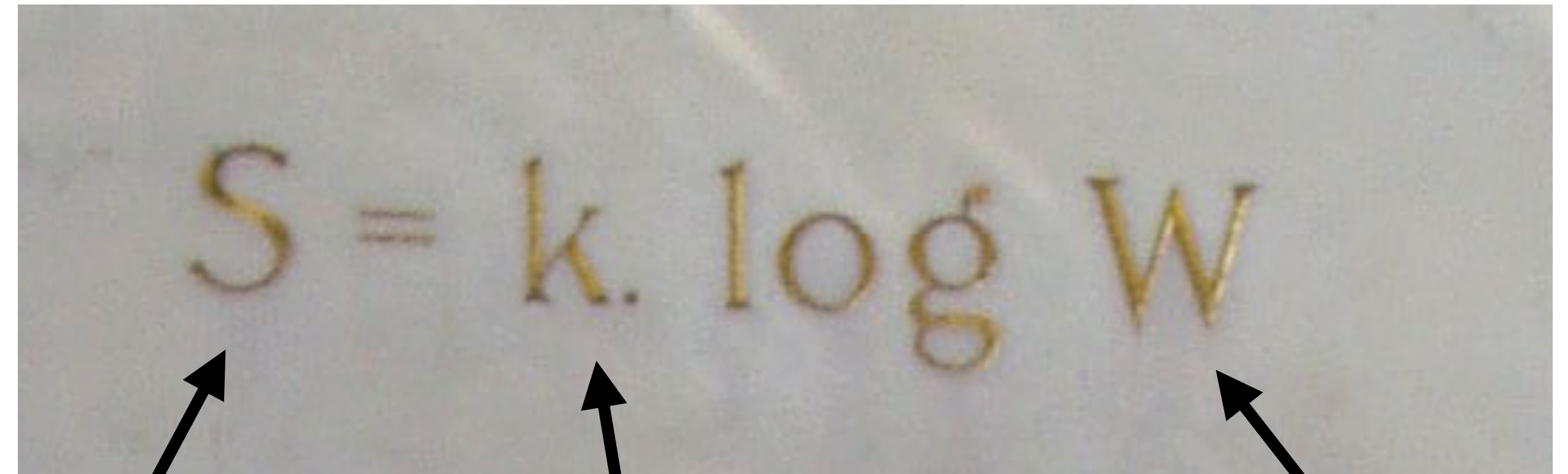
Most likely values  $\lambda$  are those for which  $\beta \tilde{H}_\lambda - \log \mu(\mathcal{O}(\mathbf{X}) = \lambda)$  is as **small** as possible

*Free energy*: “ $\beta \times$  Energy – logarithmic volume (Entropy)”

The **minimization** problem clearly **depends on  $\beta$**  (the inverse temperature)

Does it depend on  $\beta$  in a “dramatic” way? *Phase transition?*

# Entropy = logarithmic volume



**Entropy**

**Boltzmann's  
constant**

**Macrostate  
volume**

*Wahrscheinlichkeit*

# Models

# Lattice models

- **Ising model** on a lattice  $\Lambda$  (+ many related models)  
State space =  $\{-1, +1\}^\Lambda$  “spins”  
Energy = **alignement** vs. non-alignement of neighboring spins  
Observable = **magnetization**

Extremely well-studied

Non-trivial geometric questions (e.g. describe the **+1/-1 regions**)  
...but not very interesting in terms of “point configurations”

# Continuum models

- **Particles**  $x_1, \dots, x_N$  in physical space (a domain of  $\mathbb{R}^d$ , a manifold...)
- Interactions between each **pair of particles** via a *potential*  $\varphi$   
$$\text{Energy} = \sum_{i \neq j} \varphi(x_i - x_j)$$
- Introduce the **Gibbs measure = law of a random point configuration in space !**  
+ Choose an observable characterizing **order / disorder**.
- Existence of *phase transitions* in the **continuum** is on B. Simon's 1984 list of  
« Fifteen problems in mathematical physics »

# Goal: phase diagram

Will depend on dimension  $d$  and on interaction  $\varphi$

$$T = 0, \beta = +\infty$$

$$T = +\infty, \beta = 0$$

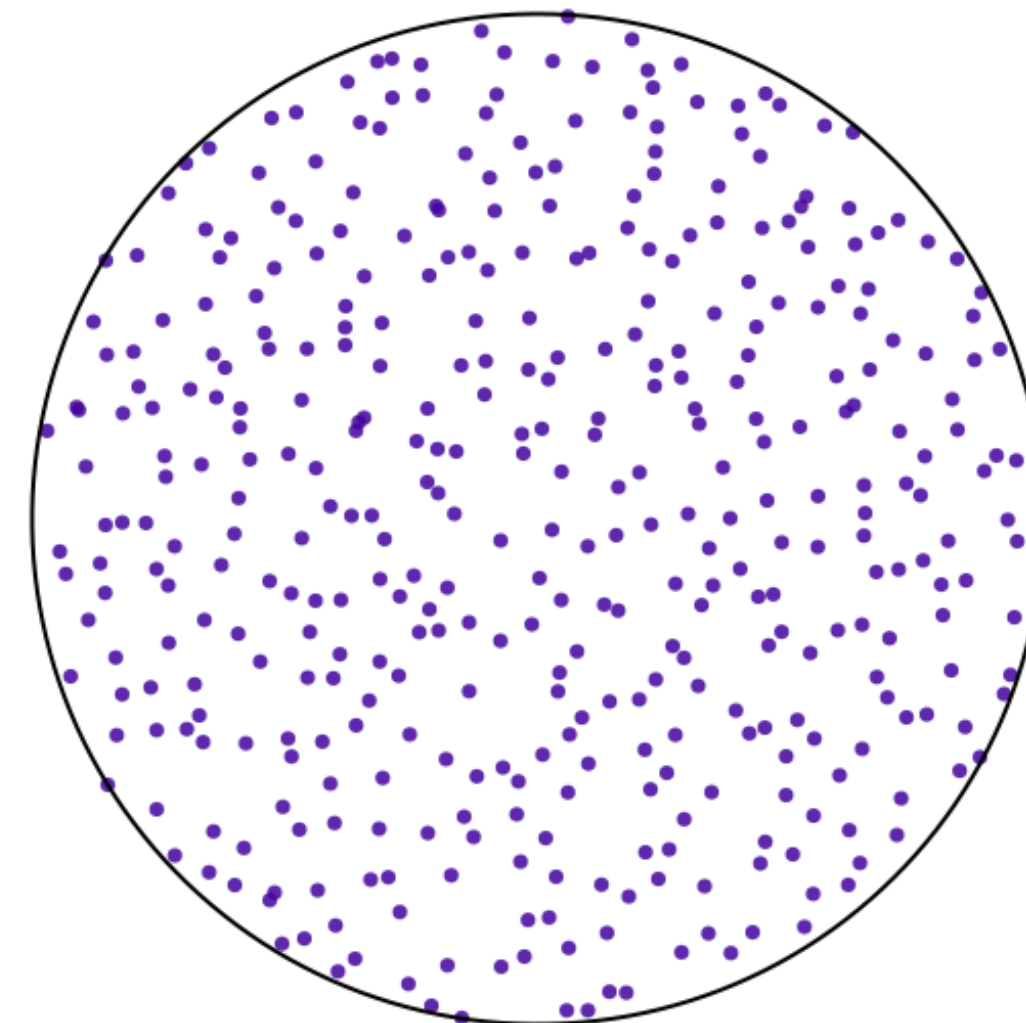
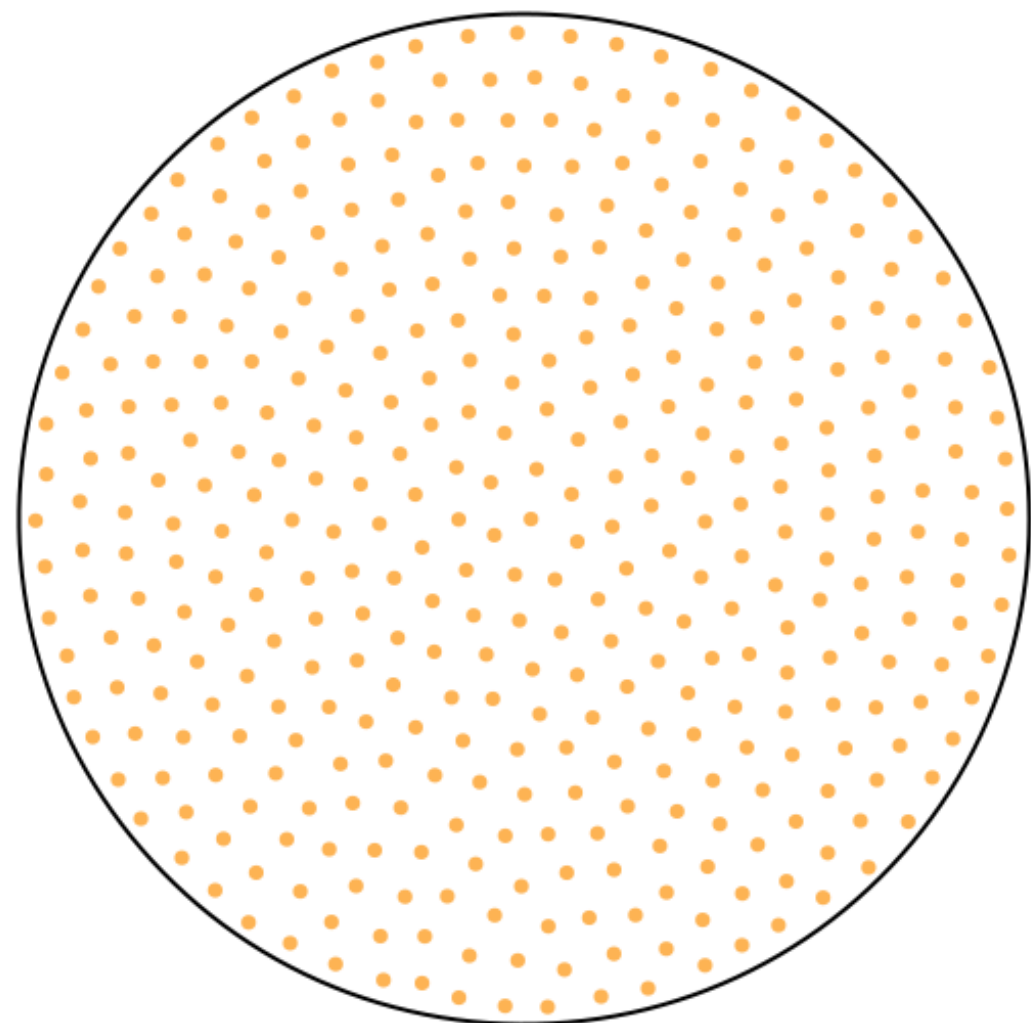
Lattice (?)

Order

Transition?

Disorder

Independent particles



$$\text{Energy} = \sum_{i \neq j} \varphi(x_i - x_j), \text{ who is } \varphi ?$$

# Riesz interactions

Power-Law  $\varphi(x) = \frac{1}{|x|^s}$ , *Riesz potential*

In dimension  $d$  :

$s > d$  short-range

$s < d$  long-range

$s = d - 2$  Coulomb

$s = 0 \leftrightarrow \varphi(x) = -\log|x|$

Expect **major differences** between short- and long-range  
e.g. when  $s < d$ , the energy ceases to be (almost-)additive w.r.t. space

# Other interactions

- Lennard-Jones  $\frac{1}{|x|^{12}} - \frac{1}{|x|^6}$  *non-monotonous*
- Gaussian interaction (*Gaussian core model*)
- **Hardcore** interactions ( $\varphi(x) = +\infty$  if  $|x| \leq R_0$ )
- A **general interaction** + assumptions on *regularity and range* (60's...)

$$d\mathbb{P}_{N,\beta}(x_1, \dots, x_N) = \frac{1}{Z_{N,\beta}} e^{-\beta \sum_{i < j} \varphi(|x_i - x_j|)} d\mu(x_1) \dots d\mu(x_N)$$

**Finite volume**

# Gaussian $\beta$ -ensemble

**RMT origin:** take a  $N \times N$  **Hermitian** matrix with **Gaussian coefficients**  
**Joint distribution** of eigenvalues is **computable** Dyson 60's

→ Gibbs measure with **logarithmic** interaction... **on the line**...at  $\beta = 2$

$$d\mathbb{P}_{N,\beta}(x_1, \dots, x_N) = \frac{1}{Z_{N,\beta}} e^{-\beta \sum_{i < j} -\log|x_i - x_j|} d\mu(x_1) \dots d\mu(x_N)$$

$\mu$  is some Gaussian measure on the real line

Existence of a random matrix model for all  $\beta > 0$  (**Dumitriu-Edelman '00**)

Called  $\beta$ -ensemble,  $G\beta E$  or *one-dimensional log-gas*

# Circular $\beta$ -ensemble

Take a **uniform**  $N \times N$  **unitary** matrix

**Joint distribution** of eigenvalues is **computable**

→ Gibbs measure with **logarithmic** interaction... **on the circle**...at  $\beta = 2$

Again: RMT model for general  $\beta$

$$d\mathbb{P}_{N,\beta}(x_1, \dots, x_N) = \frac{1}{Z_{N,\beta}} e^{-\beta \sum_{i < j} -\log|x_i - x_j|} d\mu(x_1) \dots d\mu(x_N)$$

$\mu$  uniform measure on the unit circle

Called  $C\beta E$  or *one-dimensional periodic log-gas*

# Two-dimensional log-gas

Take a  $N \times N$  matrix with **Gaussian coefficients (no symmetry)**

**Joint distribution** of eigenvalues is **computable**

→ Gibbs measure with **logarithmic** interaction... in the plane, at  $\beta = 2$

$$d\mathbb{P}_{N,\beta}(x_1, \dots, x_N) = \frac{1}{Z_{N,\beta}} e^{-\beta \sum_{i < j} -\log|x_i - x_j|} d\mu(x_1) \dots d\mu(x_N)$$

$\mu$  is some Gaussian measure on the plane

General  $\beta$  ? No RMT interpretation, but **existence in statistical physics.**

Called (one-component) **Coulomb gas** or **one-component plasma**

# Quick summary

- Gaussian  $\beta$ -ensemble / **1d log-gas**: finite random point configuration on the **line**
- Circular  $\beta$ -ensemble: finite random point configuration on the unit **circle**
- **2d log-gas**: finite random point configuration in the **plane**

(Those three have logarithmic interaction, and some link with RMT)

- **Riesz gases** in arbitrary dimension: **short/long range**, **Coulomb case**  $s = d - 2$
- General potential  $\varphi$ : finite random point configuration in **dimension**  $d$

# Some friends

2d log-gas

= Particles + logarithmic interaction in  $\mathbb{R}^2$

= Random *complex* eigenvalues ( $\beta = 2$ )

1d log-gas

= Particles + logarithmic interaction in  $\mathbb{R}$

= Random *real* eigenvalues ( $\beta > 0$ )

## Random polynomials and Gaussian entire functions

Take  $(a_k)_{k \geq 0}$  independent complex Gaussian coefficients, and form:

$$f(z) := \sum_{k \geq 0} \frac{a_k}{\sqrt{k!}} z^k$$

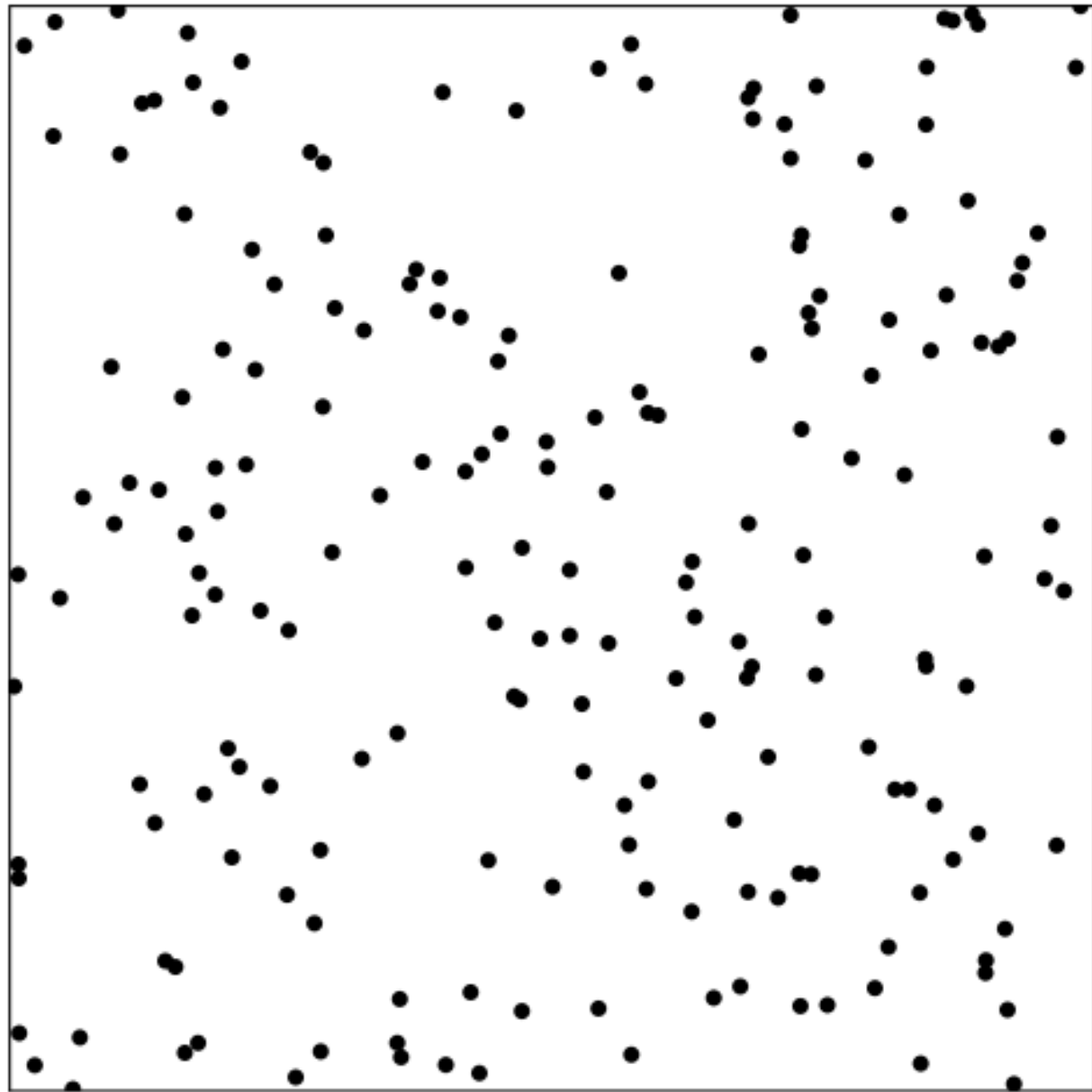
*Gaussian Entire Function (GEF)*

Zeros of  $f$  share many similarities  
with a 2d log-gas

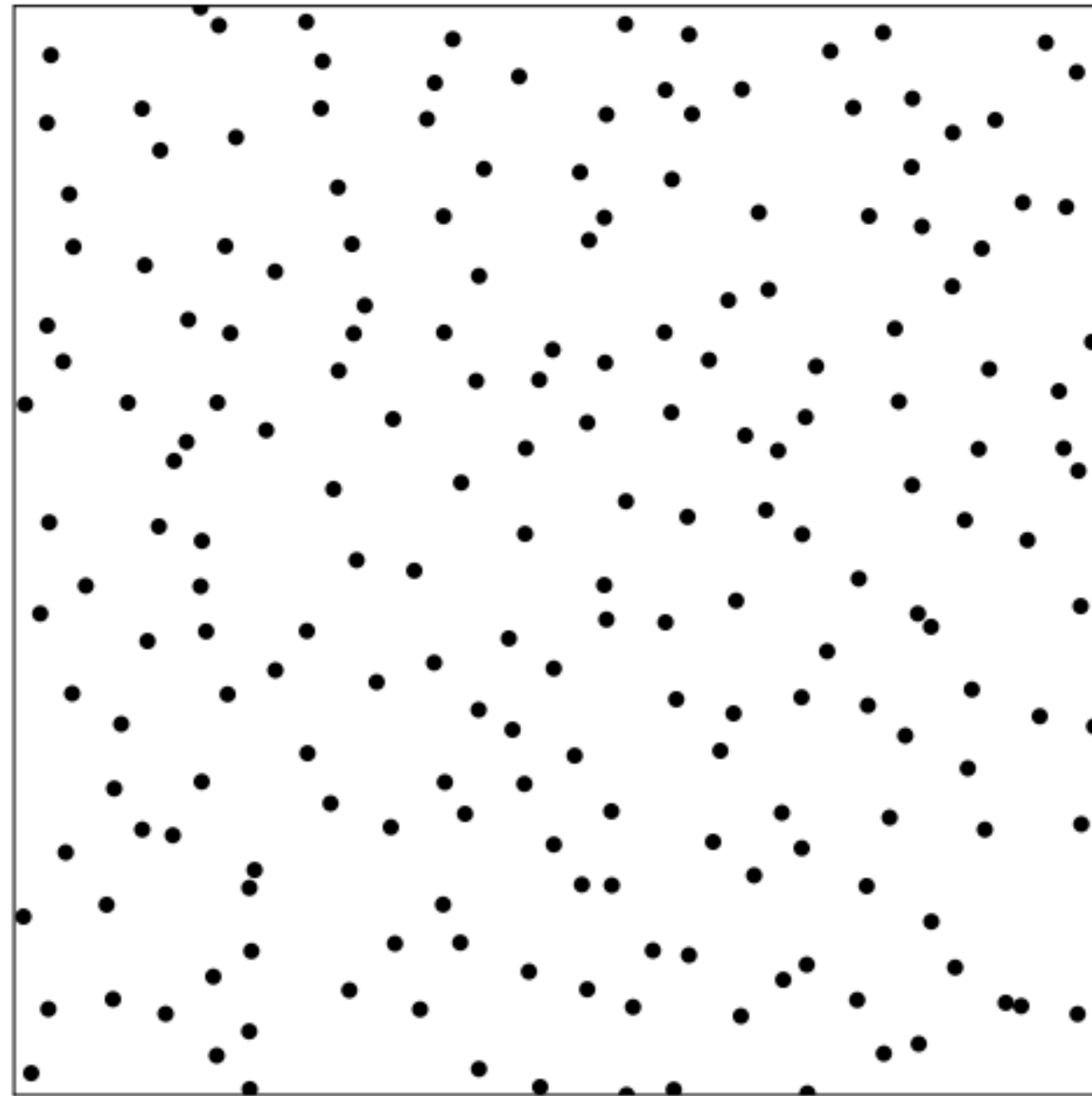
$$P_N(z) := \sum_{k=0}^N a_k z^k$$

*Gaussian Kac Polynomial*

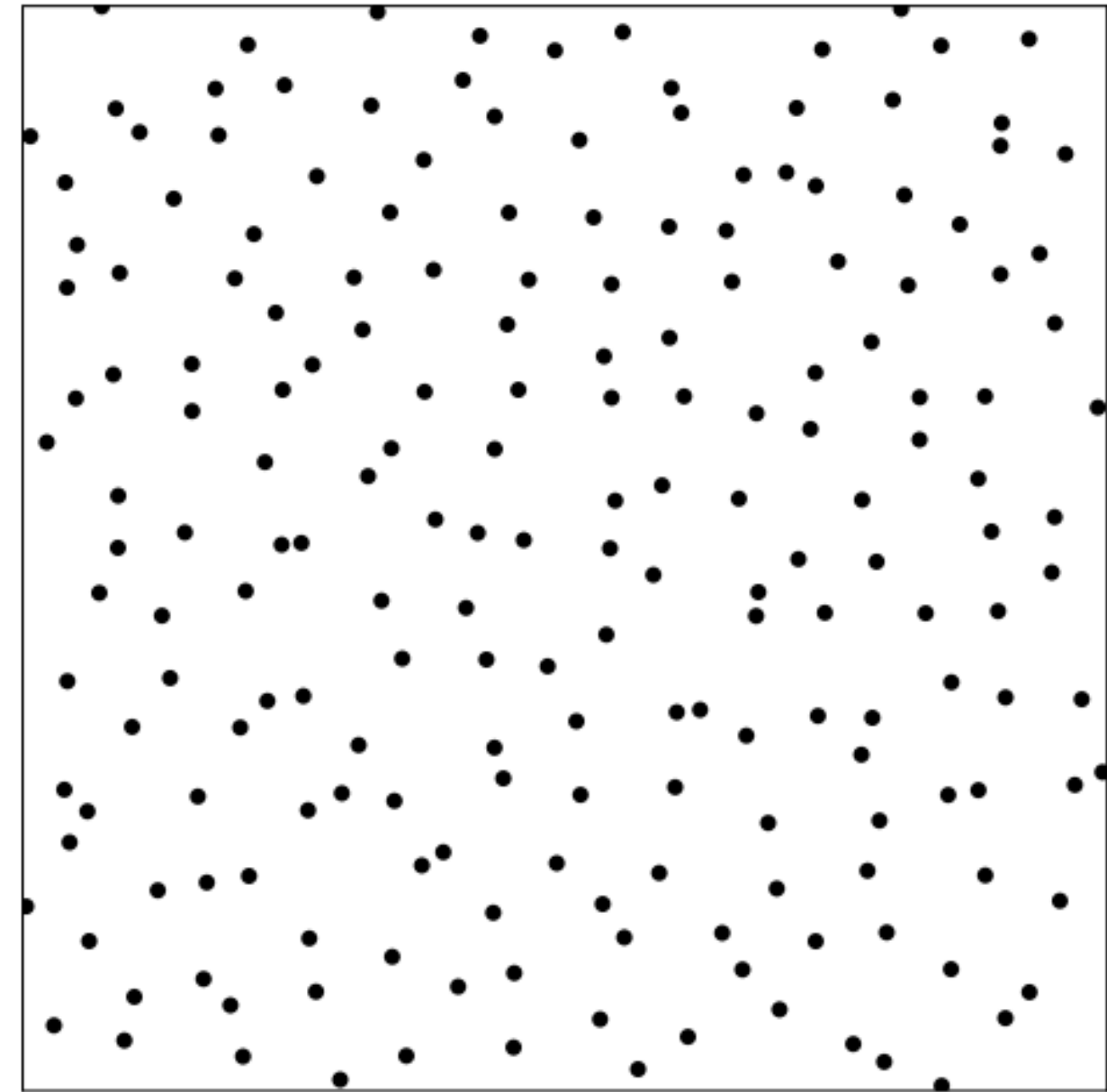
Zeros of  $P_N$  share many similarities  
with a 1d log-gas



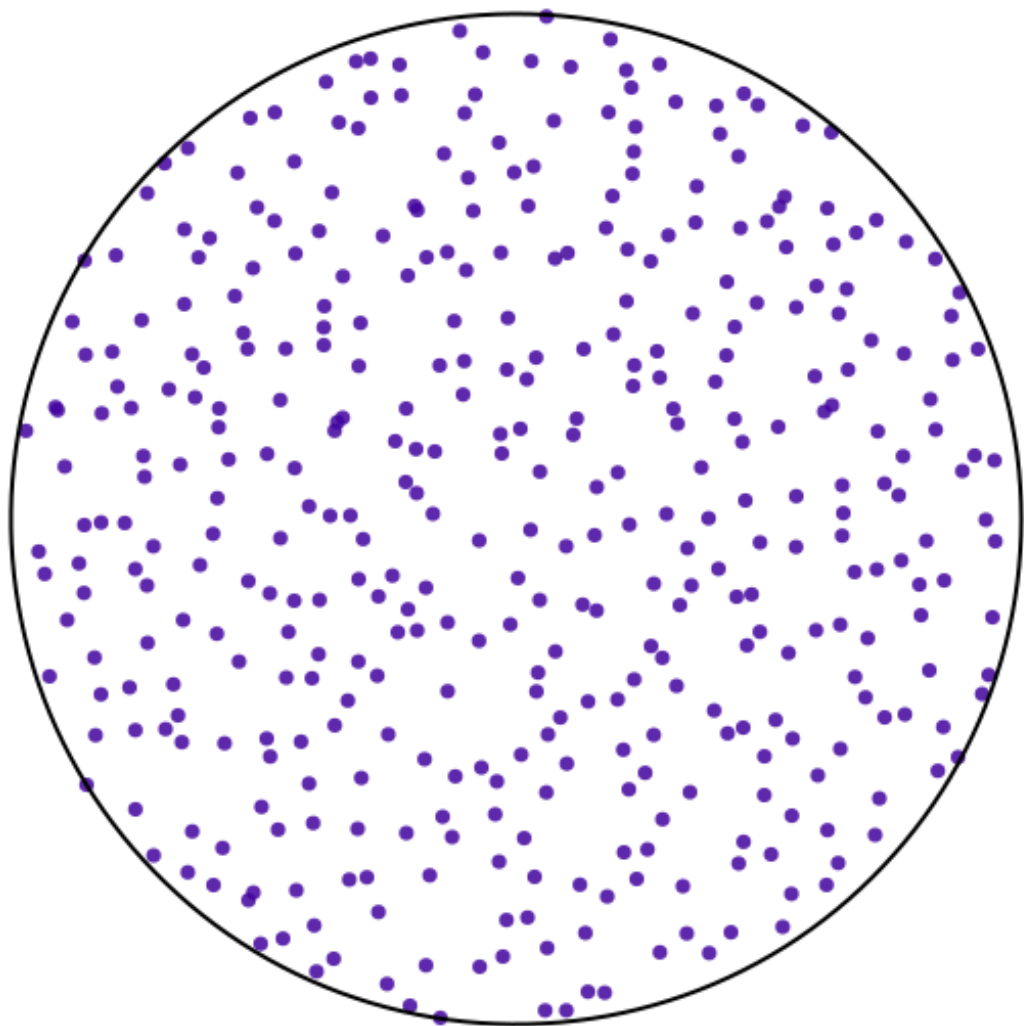
Independent points



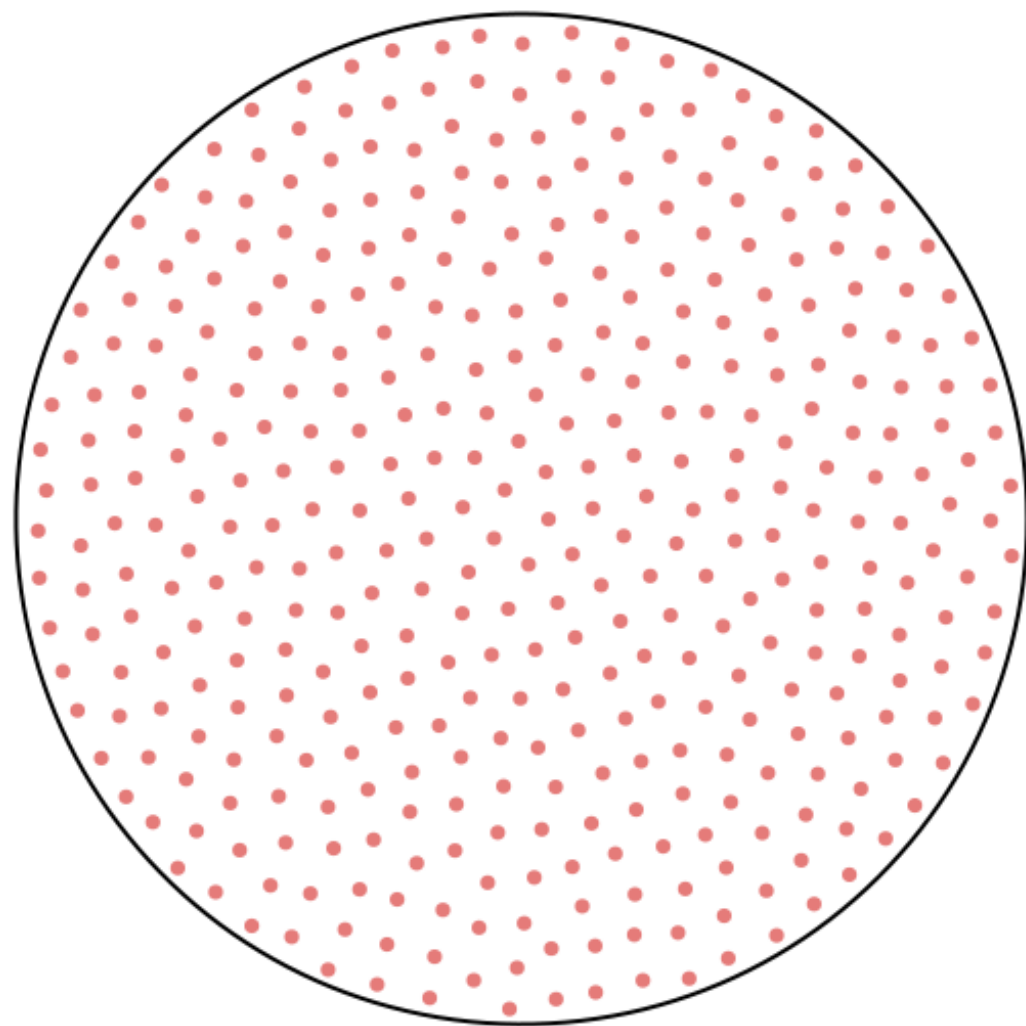
Ginibre ensemble (2DOCP)



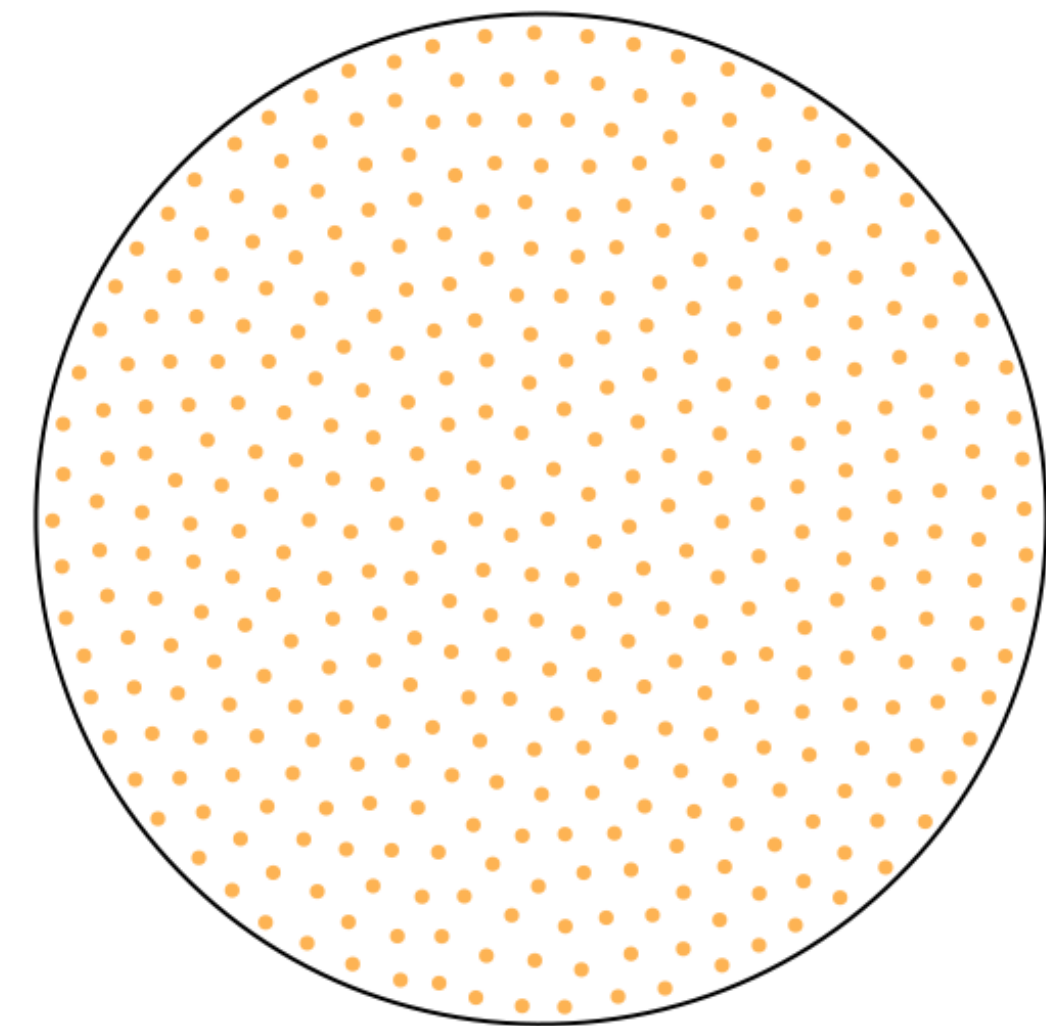
Zeroes of GEF



$\beta$  small



A 2d log-gas



$\beta$  large

# Infinite-volume limit

- We would like to let  $N \rightarrow \infty$ , obtain **infinite** random point configuration
- Existence of a limit / limit points is **hard in general!** *Uniqueness matters.*
- Known:
  - For  $\beta$ -ensembles (Gaussian and Circular): **Sine- $\beta$  point process** (Valkó-Virág '09)
  - For 2d log-gas at  $\beta = 2$ : **Ginibre point process** ('64), **general  $\beta$  is open**
  - For some long-range 1d Riesz gases: **Riesz- $\beta$  point process** (Boursier '21)

Instead of studying the Gibbs measure  $\mathbb{P}_{N,\beta}$  as  $N \rightarrow \infty$ , study the *infinite-volume variational problem of minimizing free energy?*

# Free energy minimization


Dimension  $d$

Config = point configurations in  $\mathbb{R}^d$

$\mathcal{P}(\text{Config}) =$  probability measures on Config

Given the interaction potential  $\varphi$ , **define** (??) an *energy density*  $H_\varphi : \text{Config} \rightarrow \mathbb{R}$

Minimize  $\beta \mathbb{E}_P[H_\varphi] + \mathcal{E}(P)$  over  $\mathcal{P}(\text{Config})$



Average energy density under  $P$

Entropy density of  $P$

In classical cases, this describes the infinite-volume limit “equilibrium” state(s)

# Random point configurations

This leads us to study *random point configurations* in some generality:

- Because one can hope to solve the **free energy minimization** problem
- To derive good notions of **order** and **disorder**
- Because they might be *good candidates for deterministic questions!*
- Because it's **fun**? *Stochastic geometry.*

# Some classes and operations

- Determinantal point processes (DPP): correlation functions expressed through a *determinant*  $\rho_n(x_1, \dots, x_n) = \det(K(x_i, x_j))$  *Explicit computations!*  
Examples: 1d log-gas at  $\beta = 2$ , 2d log-gas at  $\beta = 2$
- **Perturbed lattices**: move lattice points at random. *Large class!*
- **Gibbs p.p.**: have a local “density” of the form  $e^{-\beta H}$  for some *energy*  $H$
- Operations: superposition / thinning, mixture. **Displacement interpolation?**

Order / Disorder for random point configurations

# Order through correlations

1-point correlation  $\rho_1(x)$  = “probability” of having a particle at  $x$

2-point  $\rho_2(x, y)$  = “probability” of having a particle at  $x$  **and** a particle at  $y$

**Decay of correlations ?**

$$\rho_2(x, y) - \rho_1(x)\rho_1(y) \rightarrow 0 \text{ as } |x - y| \rightarrow \infty$$

**Exponential** (« fast ») decay = “**disorder**”

**Polynomial** (« slow ») decay = “quasi-order”

**No decay** = “**order**”

**Repulsion ?**

$$\rho_2(x, y) \ll \rho_1(x)\rho_1(y) \text{ for } |x - y| \ll 1$$

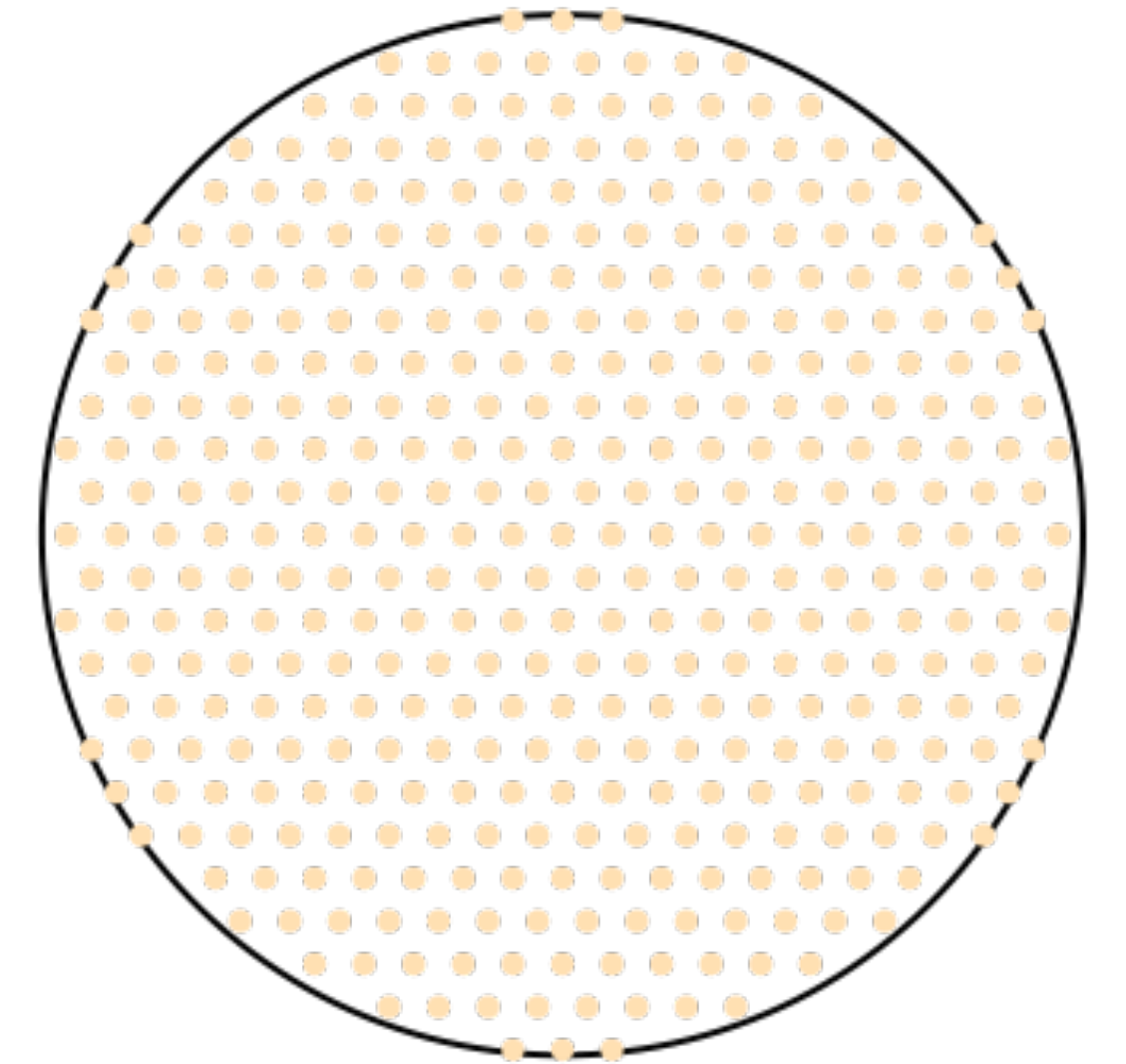
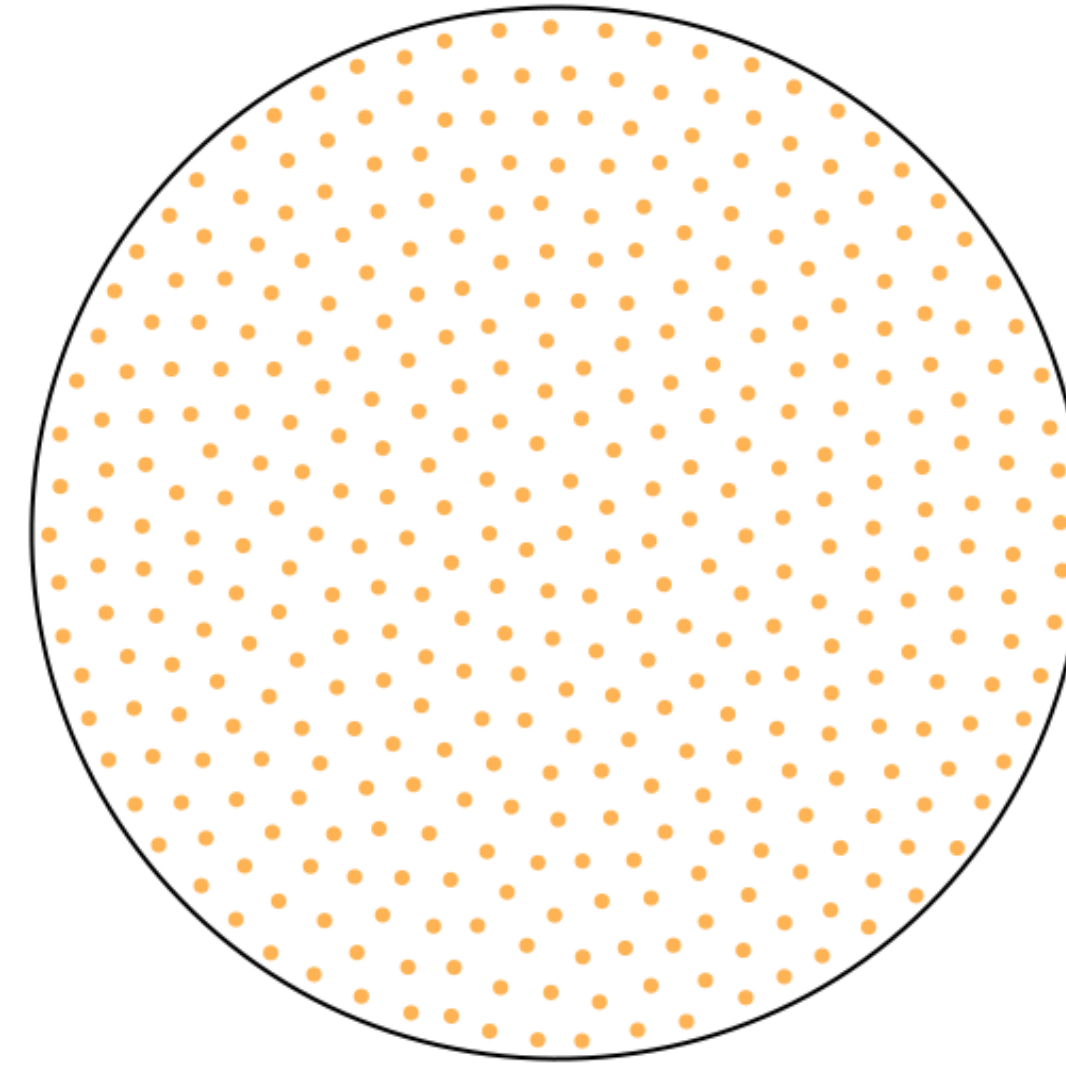
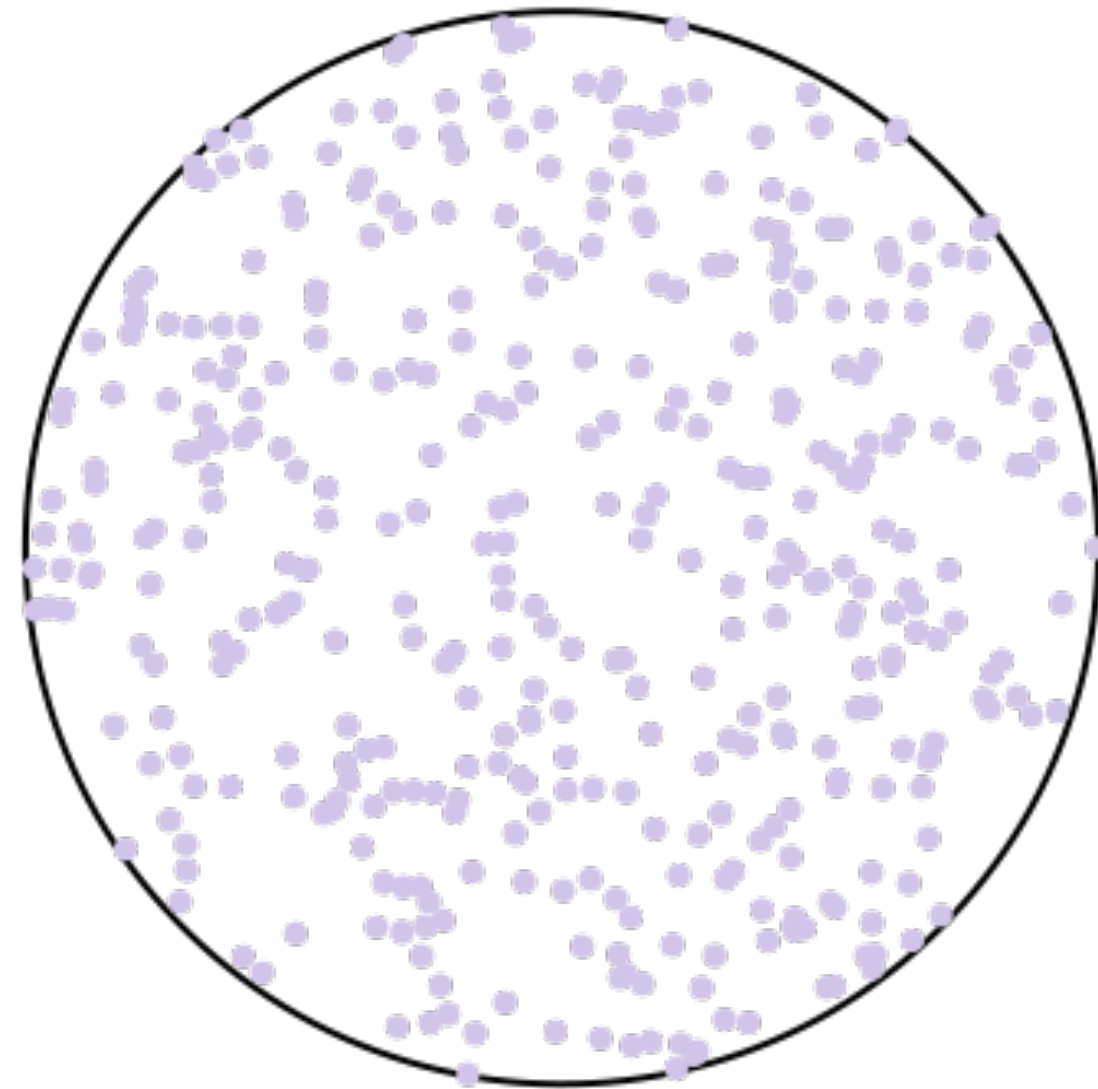
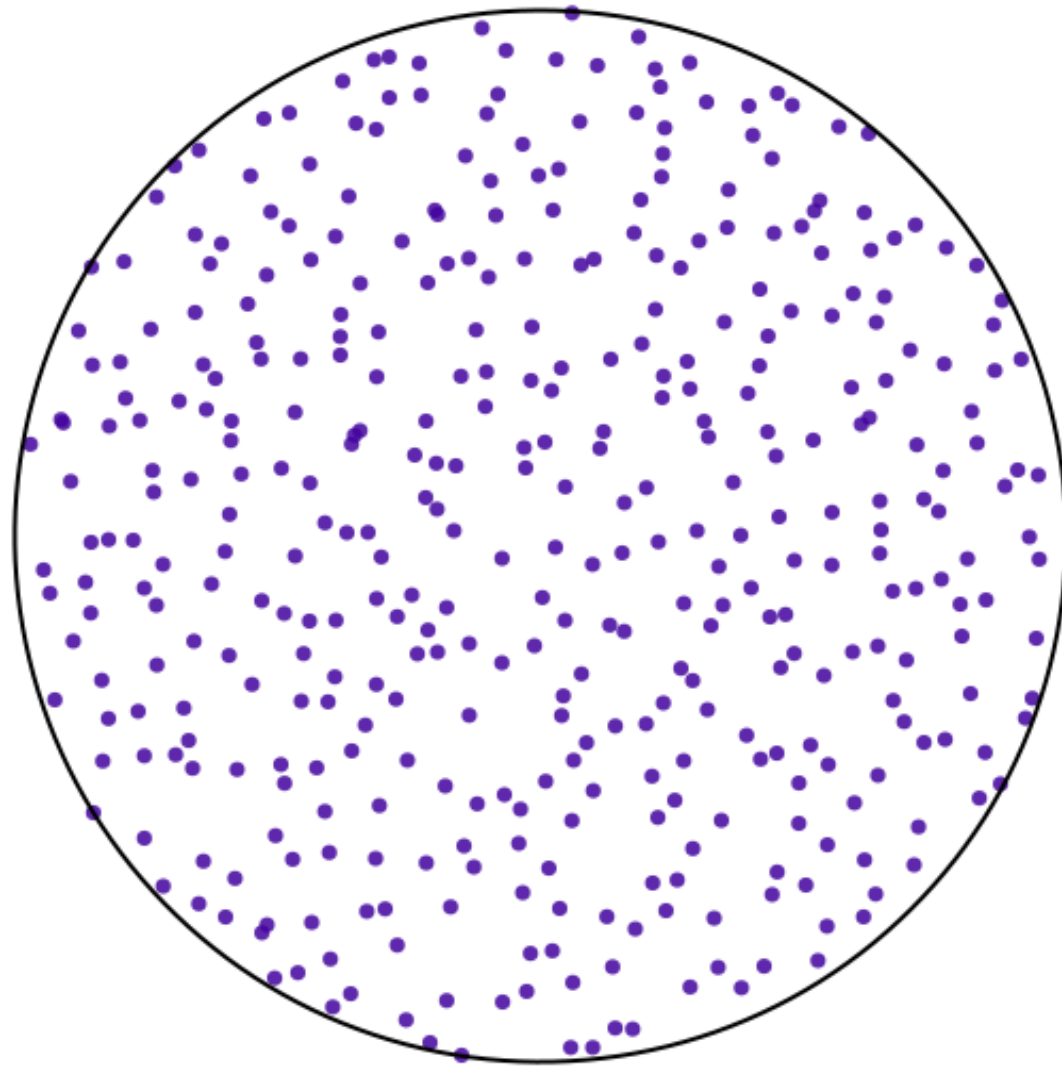
Determinantal point processes are repulsive

# But: correlations are hard

- **Correlation functions** - or simply their decay - are *very hard to compute*
- **Recent progress for the 1d log-gas**
- For the 2d log-gas, only known case for 2-point correlation is  $\beta = 2$ :
  - decay is  $e^{-|x-y|^2}$  for  $|x - y| \gg 1$  (very fast!)
  - vanishes like  $|x - y|^2$  for  $|x - y| \ll 1$  (*repulsion*)

Moreover, some regularity of those systems is **not** captured  
by the usual “decay of correlations” criterion  
One sometimes speaks of *order within disorder*

# Other faces of order



1. Size of fluctuations
2. Rigidity with respect to the exterior
3. Transportation cost

# Size of *fluctuations*

Take a big ball of radius  $R \gg 1$  and count **how many points fall into it**.

What is the *variance* of this quantity (measures “*charge fluctuations*”)?

For **independent** points: grows like  $R^d$ .

For a **lattice** with independent perturbations, or uniform shift: like  $R^{d-1}$ .

Physics literature: **Anything with variance  $o(R^d)$  is interesting.**

*Hyperuniform* or *super-homogenous* systems. **Torquato.**

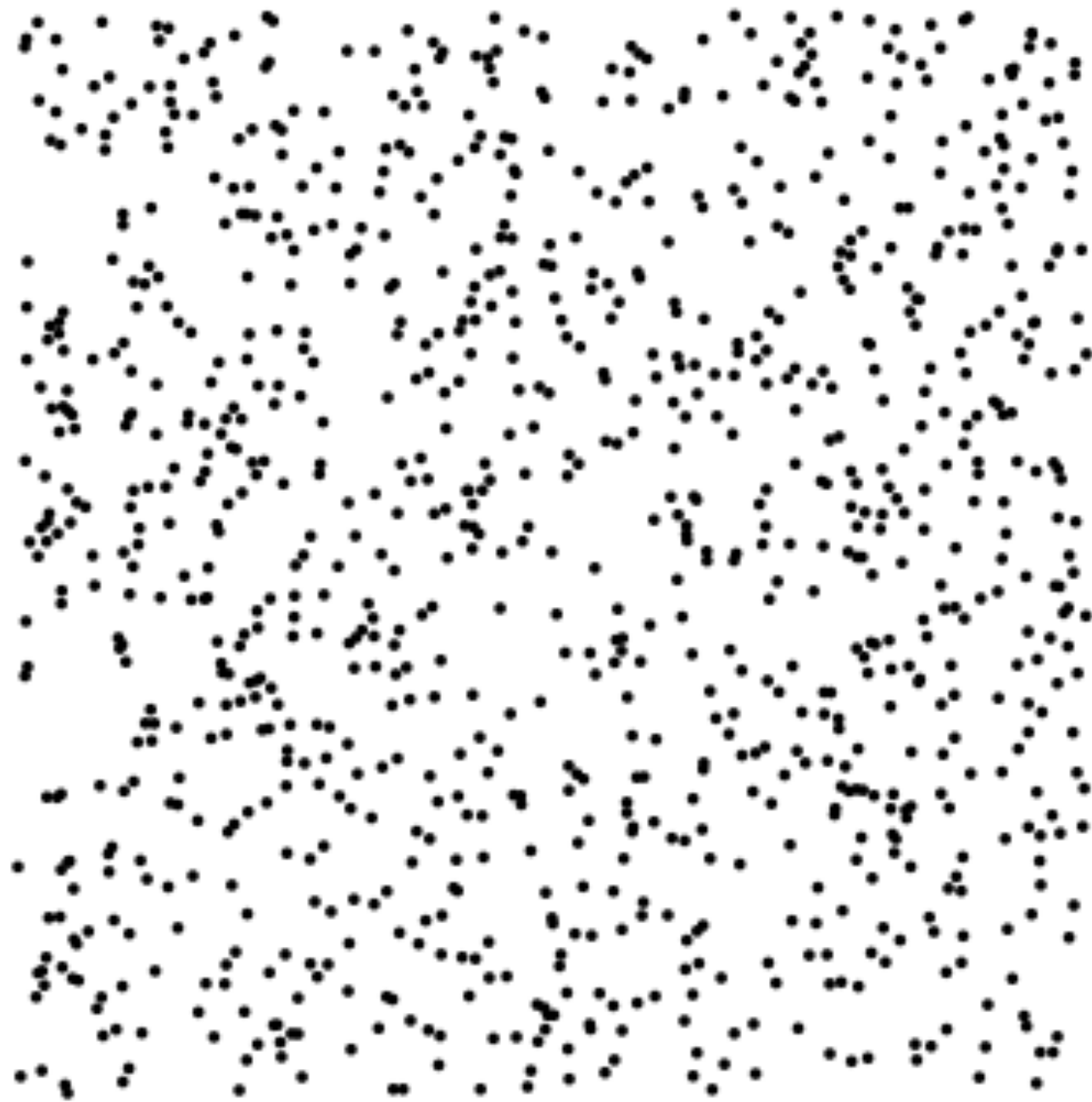
Instead of counting points, what about

Variance  $\left[ \sum_{i=1}^N f(x_i) \right]$  for a nice function  $f : \mathbb{R}^d \rightarrow \mathbb{R}$ ?

**Fluctuations of linear statistics**

# Rigidity properties

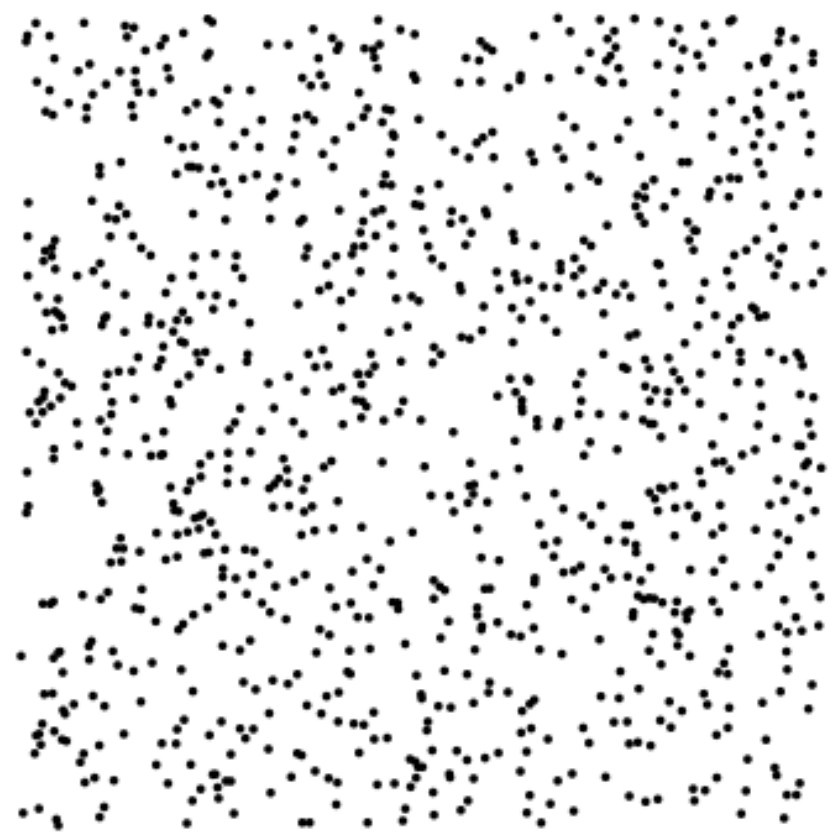
Ghosh-Peres '12



What can we say about  
the *inside* region  
from looking *outside*?

- (Almost surely) guess the **number** of points? *Number-rigid.*
- (Almost surely) guess the **center of mass**? *Barycenter-rigid.*
- (Almost surely) guess the full **configuration**? *Fully-rigid.*

# Who is rigid?



- A Poisson point process (independent points)  
*No rigidity*
- A lattice (perfectly ordered)  
*Fully-rigid*

What about lattice + i.i.d. perturbations?

Surprisingly subtle question

Peres-Sly '14

1d Log-gas is Chhaibi-Najnudel '15

**Number-rigid**  $\forall \beta > 0$

Zeros of GEF are

**number-rigid and** Ghosh-Peres '12

**barycenter-rigid**

Some 1d Riesz gases

**are not number-rigid**

Dereudre-Vasseur '18

Some are

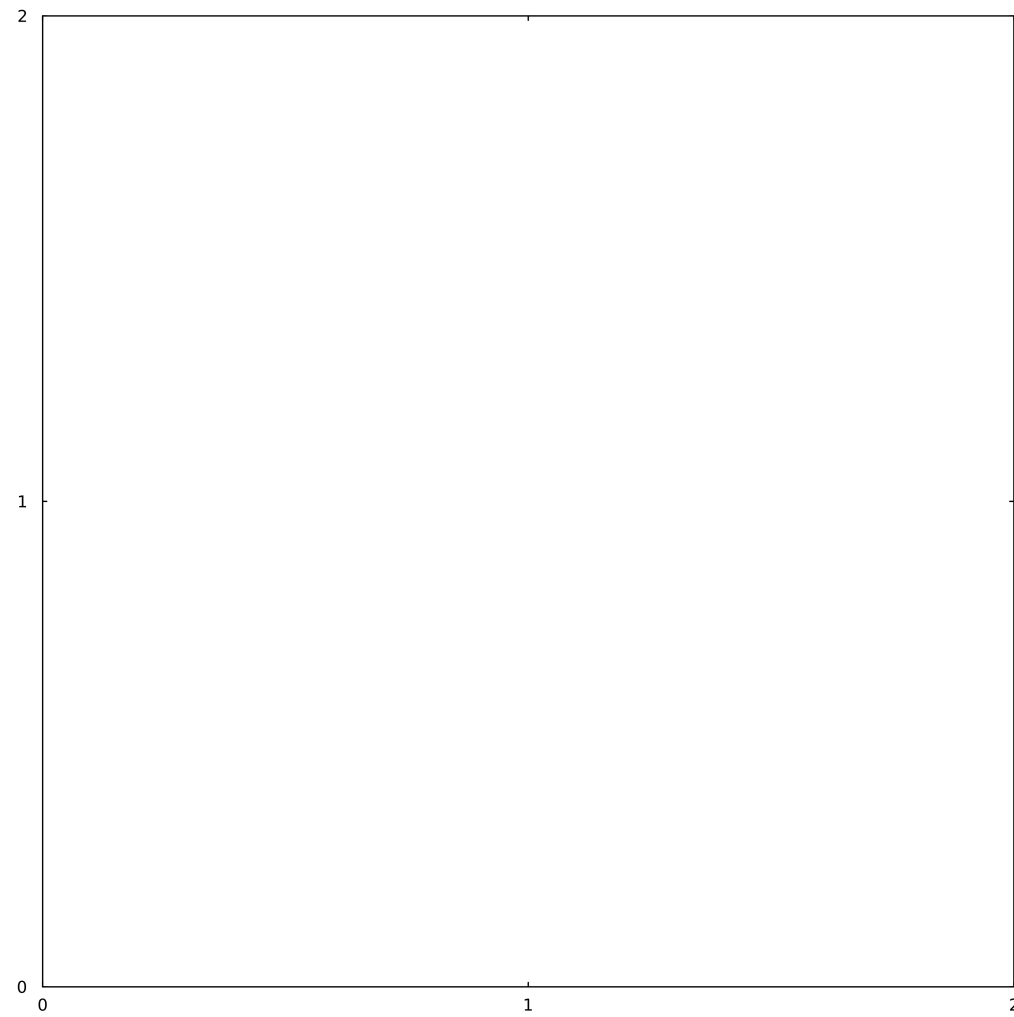
Dereudre-Digneaux '25

2d Log-gas is

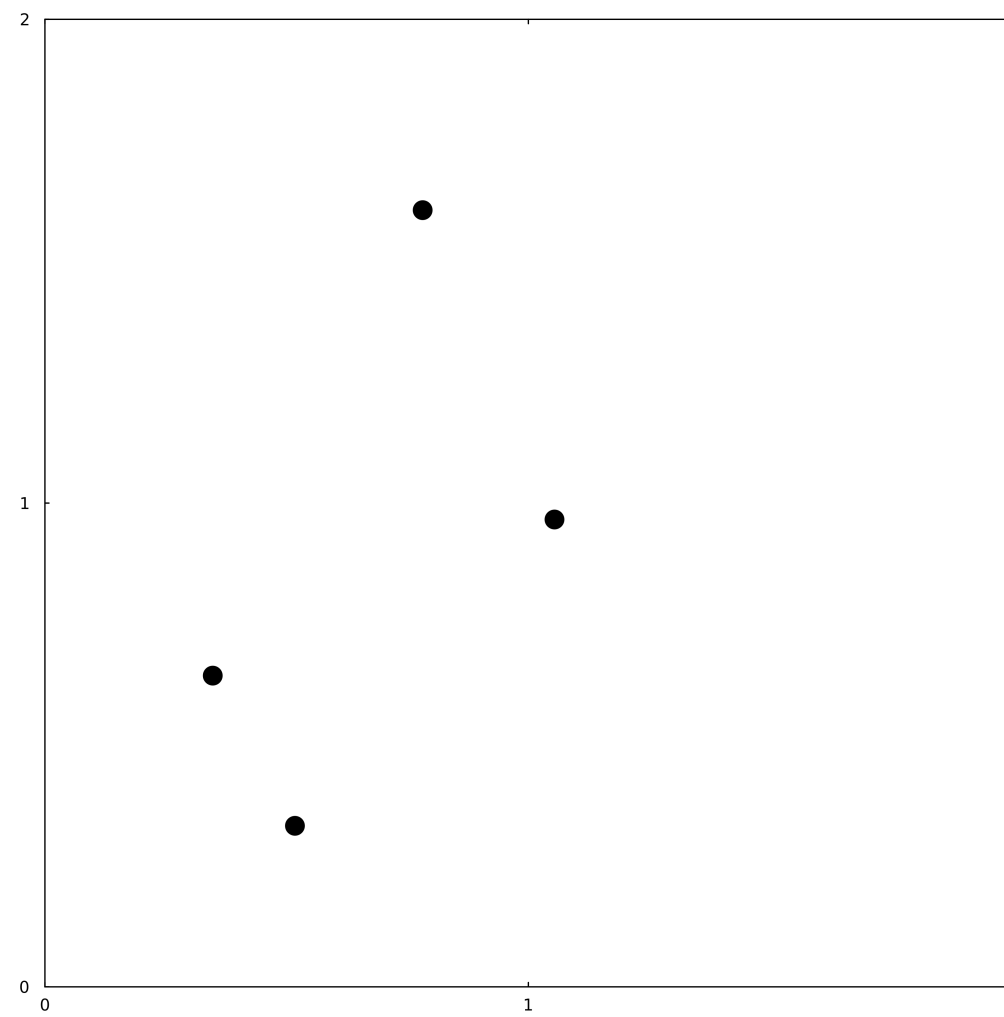
**Number-rigid**  $\forall \beta > 0$

L. '24

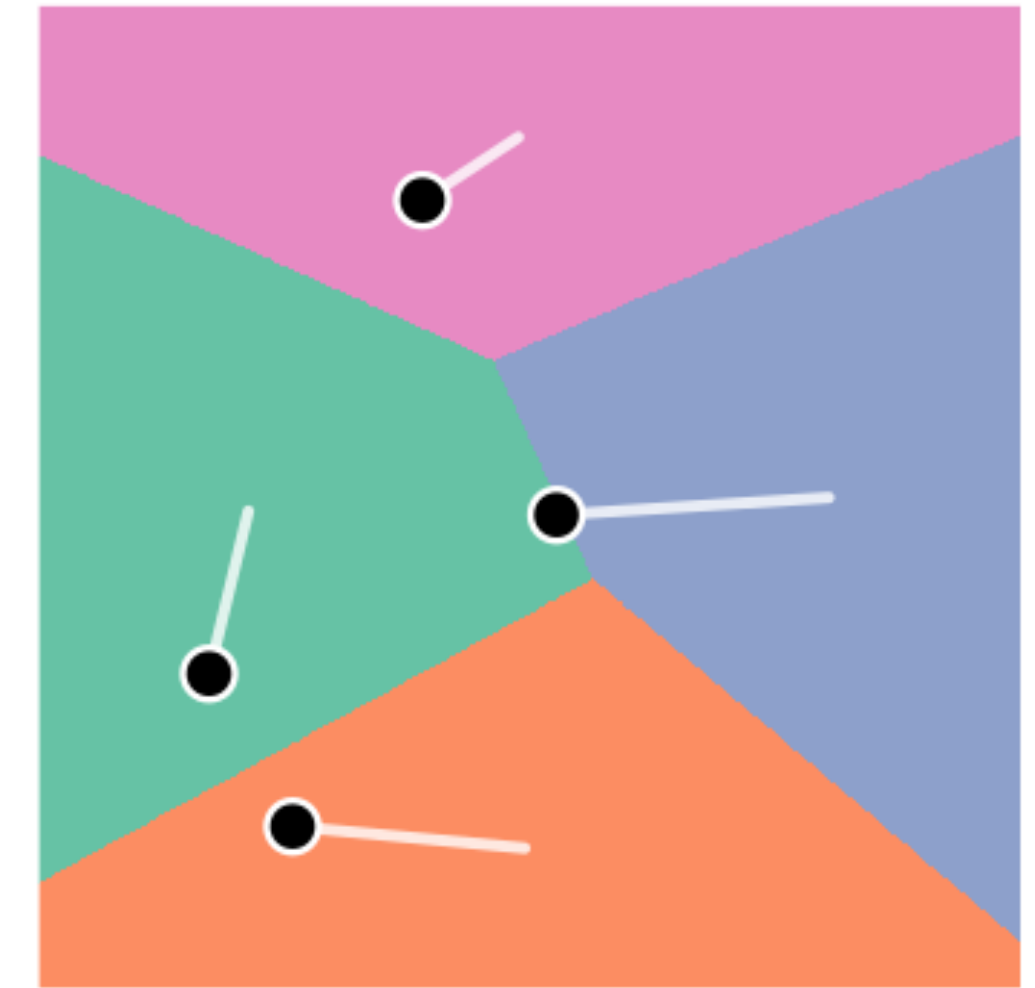
# Transportation cost



Take a  $L \times L$  square



Place  $L^2$  points



Associate to each point  $x$   
a cell  $C_x$  of area 1

*Quadratic transportation cost*  
for the  $x$ -cell:  $\int_{C_x} |x - t|^2 dt$

*Transportation cost*  
 $\sum_x \int_{C_x} |x - t|^2 dt$

What is the *minimal* cost?

# Wasserstein “order” in $d = 1, 2$

Ajtai-Komlos-Tusnady '84

If the points are on a grid (*order*)  
the cost (per point) is **bounded**

If points are drawn i.i.d. at random (*disorder*)  
average cost **diverges as  $L \rightarrow +\infty$**  in  $d = 1, 2$ .

In dimension **3 and above**, this does not  
capture a meaningful difference between  
order and disorder

# Summary of order / disorder

- “Classical” notion of order through correlation functions
- Some alternative notions: hyperuniformity, small fluctuations of linear statistics, rigidity *à la* Ghosh-Peres, Wasserstein distance to Lebesgue
- **Model-dependent** answers but few (if any)  $\beta$ -dependent answers yet.

# Final word: the $\beta = +\infty$ case

Minimizing “ $\beta \times \text{Energy} - \text{logarithmic volume (Entropy)}$ ”

For  $\beta = +\infty$ , this boils down an *energy minimization problem*

*Deterministic, but notoriously difficult...*

Hard to study  $\beta \gg 1$  when you don't know  $\beta = +\infty$ ...

*Free energy minimization is an “extension” of this problem in a random setting.*

Thank you for your attention