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- Webpage for OOPS: https://www.math.ubc.ca/Links/OOPS/.
- Source is available at https://github.com/holdenlee/oops. Contributions, corrections, and comments welcome; feel free to send a pull request. Any errors are probably due to me! You can ping me on zulip or email me at holden.lee@duke.edu.

A direct link to these notes is https://www.dropbox.com/s/u4lpayi98r06un5/oops.pdf?dl=0.

1 Disordered systems, rank-one matrix estimation and Hamilton-Jacobi equations (Jean-Christophe Mourrat)

We consider the problem of estimating a large rank-one matrix, given noisy observations. This inference problem is known to have a phase transition, in the sense that the partial recovery of the original matrix is only possible if the signal-to-noise ratio exceeds a (non-zero) value. We will present a new proof of this fact based on the study of a Hamilton-Jacobi equation. This alternative argument allows to obtain better rates of convergence, and also seems more amenable to extensions to other models such as spin glasses.

References:

- Paper: [Mou18], https://cims.nyu.edu/~jcm777/HJinfer.pdf
 Approach based on weak solutions: [Mou19], https://cims.nyu.edu/~jcm777/HJrank.pdf
- Book for background material: [FV17], available at https://www.unige.ch/math/folks/velenik/smbook/index.html

• Other approaches

- Lelarge-Miolane: [LM19], https://arxiv.org/abs/1611.03888

- Barbier-Macris: [BM19], https://arxiv.org/abs/1705.02780

2020/5/18 Lecture 1

Suppose students are assigned one of two dormitories. They put on a sorting hat, which decides which dorm they go in.

The students are $i \in \{1, ..., N\}$. An assignment is $\sigma \in \{\pm 1\}^N$. The sorting had optimizes the quality of interaction between i and j, J_{ij} . Suppose (J_{ij}) are independent standard Gaussians. The larger J_{ij} is, the more that i and j like to be together. We want to maximize $\sigma \mapsto \sum J_{ij} \mathbb{1}_{\{\sigma_i = \sigma_j\}}$. By a linear transformation this is equivalent to maximizing $\sum J_{ij}\sigma_i\sigma_j$.

What is $\max_{\sigma\{\pm 1\}^N} \sum J_{ij}\sigma_i\sigma_j$ as $N \to \infty$? Because the J_{ij} can be positive or negative, we can't make all the students happy. Thus we can say there are **frustrations** in the problem. These models are called spin glasses in the literature. It's difficult to find the optimum: making local moves, you may have to decrease the objective before increasing it.

I want to consider a softer version of the maximum. We look at the spin glass model¹

$$\mathbb{E}\frac{1}{N}\log\sum_{\sigma\in\{\pm 1\}^N}\exp\left(\frac{\beta}{\sqrt{N}}\sum_{i,j=1}^N J_{ij}\sigma_i\sigma_j\right)$$

If β is large this is dominated by the maximum. This is like a relaxation of the problem. We should expect what's inside is order N, so we divide by N.²

Parisi in the late 70's (1979) proposed an answer for what this becomes as $N \to \infty$. It's a fairly complicated formula.

Guerra 03 and Talagrand 06 proved it rigorously. I find it mysterious; I want to think about a slight variation of the problem. Instead of connections between each i, j, think of them organized in two layers; there are interactions between but not within the layers (the graph is bipartite). This seems an innocent modification, but I could not understand what to write instead of the complicated formula!

Now I consider rank-one matrix esimation/inference. The question is statistical: we only observe a noisy version of a rank-one matrix. Can we recover information about it?

Here are some concrete settings where this is useful:

• You are Netflix, you want to make recommendations for your customers. A simple model is that whether or not a person likes a movie is captured by a few parameters of the movie (action, introspection, sad/happy, etc.) and customer, and is a linear

Note that the normalization is $\frac{1}{N}$ instead of $\frac{1}{\sqrt{N}}$ because here the $J_{ij}=1$.

²Can we fix the number of +1's and -1's? You can change the reference measure; this can be encoded as changing the reference measure. $\beta = 0$ is summing over reference measure. $\beta \to \infty$ recovers the maximum. β small is high temperature, β large is small temperature.

function of the parameters. Then you have a large low-rank matrix. A simplification is that it's a matrix of rank 1. I'll describe rank 1, but it's not hard to generalize.

• Community detection: The US is polarized, and there is a binary variable that will predict whether two people will be friends.

The common thread is the relation with certain partial differential equations, called **Hamilton-Jacobi equations**.

The Curie-Weiss model is a simple model that can be solved in many ways. I want to emphasize the method that uses intuition with Hamilton-Jacobi equations. Next when we turn to rank-1 matrix estimation, the proof will be almost the same.

Our derivation is not standard; if you want to see a more standard derivation see Friedli and Velenik [FV17].

Can you meaningfully recover information about the rank-1 matrix? In the Ising model there is a phase transition between an ordered and disordered state. In this inference problem there is also a phase transition. When signal-to-noise ratio is too small (weak), you cannot recover meaningful information. After the threshold, you can recover partial information.

1.1 Definitions

We want to study the probability measure that to each $\sigma \in \{\pm 1\}^N$, associates a weight proportional to

$$\exp\left(\frac{t}{N}\sum_{i,j=1}^{N}\sigma_{i}\sigma_{j}+h\sum_{i=1}^{N}\sigma_{i}\right).$$

The second term doesn't have interaction; it "tilts" the σ_i , giving each a preference. Here t > 0 but $h \in \mathbb{R}$. Define the expected value

$$\langle f(\sigma) \rangle_{k,h} := \frac{\sum_{\sigma} f(\sigma) \exp\left(\frac{t}{N} \sum_{i,j=1}^{N} \sigma_{i} \sigma_{j} + h \sum_{i=1}^{N} \sigma_{i}\right)}{\sum_{\sigma} \exp\left(\frac{t}{N} \sum_{i,j=1}^{N} \sigma_{i} \sigma_{j} + h \sum_{i=1}^{N} \sigma_{i}\right)}.$$

The subscripts are omitted when clear.

Define the free energy

$$F_N(t,h) = \frac{1}{N} \log \sum_{\sigma} \exp \left(\frac{t}{N} \sum_{i,j=1}^{N} \sigma_i \sigma_j + h \sum_{i=1}^{N} \sigma_i \right)$$

You might say: this is the normalization constant, we care about the measure. This is misleading because the normalization constant is the generating function of quantities we care about. You are calculating the exponential (moment) generating function of these variables. If you understand the mgf, you understand these quantities.

Moment generating function. Differentiating gives

$$\partial_h F_N = \frac{1}{N} \frac{\sum_{\sigma} \left(\sum_{i=1}^N \sigma_i\right) \exp\left(\frac{t}{N} \sum_{i,j=1}^N \sigma_i \sigma_j + h \sum_{i=1}^N \sigma_i\right)}{\sum_{\sigma} \exp\left(\frac{t}{N} \sum_{i,j=1}^N \sigma_i \sigma_j + h \sum_{i=1}^N \sigma_i\right)} = \frac{1}{N} \left\langle \sum_i \sigma_i \right\rangle \tag{1}$$

$$\partial_t F_N = \frac{1}{N} \left\langle \frac{1}{N} \sum \sigma_i \sigma_j \right\rangle = \left\langle \left(\frac{1}{N} \sum \sigma_i \right)^2 \right\rangle. \tag{2}$$

This model is simple; I can rewrite $\partial_t F_N$ in a simple way. The derivatives are all order 1. It's a good starting point to notice that

$$\partial_t F_N - (\partial_h F_N)^2 = \left\langle \left(\frac{1}{N} \sum \sigma_i\right)^2 \right\rangle - \left\langle \frac{1}{N} \sum \sigma_i \right\rangle^2.$$

This is the mean magnetization, the variance of the magnetization. Idea: The variance is lower-order, so as $N \to \infty$, F_N solves the equation with 0 on the right.

 F_N is the mgf of $\sum \sigma_i$. So in particular it should encode the variance of the variable in some way. I should find a way to express it in terms of F_N . Looking at the second derivative is a good idea.

$$\partial_h^2 F_N = \frac{1}{N} \left\langle \left(\sum \sigma_i \right)^2 \right\rangle - \frac{1}{N} \left(\left\langle \sum \sigma_i \right\rangle \right)^2.$$

So we have shown

$$\partial_t F_N - (\partial_h F_N)^2 = \frac{1}{N} \partial_h^2 F_N. \tag{3}$$

This is a very important observation: everything is expressed in terms of F_N . We can forget about the probability measure, definition in terms of probability measures, and just think about what F_N satisfies this equation, and what happens when N becomes large. It also suggests that as $N \to \infty$, the RHS will vanish.

I think of this as an evolution equation; think of t as time. It will be useful to understand what happens when t = 0, the initial conditions.

$$F_N(0,h) = \frac{1}{N} \log \sum_{\sigma} \sum_{\sigma} \exp\left(h \sum_i \sigma_i\right)$$
 (4)

$$= \frac{1}{N} \log \sum_{\sigma} \prod_{i=1}^{N} \exp(h\sigma_i)$$
 (5)

$$= \frac{1}{N} \log \left(e^h + e^{-h} \right)^N = \log(e^h + e^{-h}) \tag{6}$$

$$F_N(0,h) = F_1(0,h) =: \psi(h). \tag{7}$$

This does not depend on N.

The most important connection is that the h-derivative is the mean magnetization (1). What do we do with (3) and (7)?

1.2 Interlude on Hamilton-Jacobi equation

Let's take a step back and think about what the equation is saying. We need to think about what it means to be a solution of

$$\partial_t f - (\partial_h f)^2 = 0. (8)$$

The first thing to look for is a C^1 function that solves the equation pointwise. What's the problem with this?

The problem is that there is a phase transition in the Ising model. When t is small nothing impressive happens. For fixed small t, $F_N(t,h)$ will be smooth.

But for larger t, if h is positive, then the mean magnetization is positive and away from 0, and if h is tiny negative, then the mean magnetization is negative and away from 0. There will be a jump in the derivative of the function; it looks like |h|. The equation is not solved pointwise at h = 0.

QA:

• If you change the measure, you can create lots of discontinuities. Considering P with bounded support on \mathbb{R} , we can consider

$$\int \exp(\cdots) dP^{\otimes N}(\sigma).$$

You can play with P to create more corners in the limit. This changes what this ψ function.

- What's the notion of convergence for solutions as $N \to \infty$? All functions are uniformly Lipschitz. By Arzela-Ascoli there are convergent subsequences. We can take uniform convergence as the topology. If you prove convergence for some topology, you can bring it to $C^{0,1}$ topology.
- HJ equation can be solved by characteristics. Is there a probabilistic interpretation of the PDE method of characteristics? I'll try to bypass it. Barbier and Macris [BM19] use different techniques: construct characteristics for finite N. The method I present is more convenient, you don't need to follow characteristics closely, just look at whether characteristics are contracting or expanding.

2020/5/19 Lecture 2

Recall that we showed $\partial_t F_N - (\partial_h F_N)^2 = \frac{1}{N} \partial_h^2 F_N$ (3), and hoped the limit object will solve the equation (8),

$$\partial_t f - (\partial_h f)^2 = 0.$$

We have to think about what it means to solve the equation. Because of phase transitions, we don't expect solutions to be C^1 . We need to lower our expectations about what being a solution means.

Every Lipschitz function is differentiable almost everywhere. What if we just ask the equation to be satisfied almost everywhere? The problem is that the solution is not unique (given the initial value at t=0, there are multiple possible solutions). Here are some examples.

- 0 is a solution.
- $(t,h) \mapsto t+h \text{ or } t-h$.
- From these 3 solutions, we can construct another solution with the "tent function,"

$$f(t,h) = \begin{cases} t+h, & -t \le h \le 0 \\ t-h, & 0 \le h \le t \\ 0, & \text{else.} \end{cases}$$

We have an infinite number of solutions by taking combinations of translations of these.

But these are not the solutions we care about! We know the solution should be convex in the h variable; the tent function is not convex. If we add in this condition, then we get a unique solution.

Definition 1.1: We say that a Lipschitz function $\mathbb{R}_+ \times \mathbb{R} \to \mathbb{R}$ is a **weak solution** of the HJ equation (8) if

- 1. (HJ) is satisfied almost everywhere.
- 2. For every $t \ge 0$, the mapping $h \mapsto f(t, h)$ is convex.³

Proposition 1.2 (Uniqueness): If f and g are two weak solutions of the HJ equation with $f(0,\cdot)=g(0,\cdot)$, then f=g.

Why do we want this? This proposition says that two solutions with same initial condition are equal. What we want is actually a refinement of the statement: we want to compare finite-N solution (to (3)) for N large to the solution of the HJ equation. Instead of two solutions being equal, we want to show the almost-solution and the solution are close.

Proof sketch. Denote w = f - g. We have that almost everywhere

$$\partial_t w = (\partial_h f)^2 - (\partial_h g)^2$$
$$= \underbrace{(\partial_h f + \partial_h g)}_{=:h} \partial_h w.$$

Then

$$\partial_t w - b \partial_h w = 0$$
 (a.e.).

³It suffices to be semi-convex: there is a lower bound on the Hessian.

Because f, g are convex in h, the derivative of b is positive.⁴ The rough idea is to look at how the integral I(t) evolves:

$$I(t) = \int w(t, h) \, dh.$$

Suppose this is well-defined. Then integrating by parts (suppose there are no boundary terms)

$$\partial_t I(t) = \int \partial_t w = \int b \partial_h w$$
$$= -\int \underbrace{\partial_h b}_{>0} w$$

This says that I(t) wants to come back to 0. Some problems with this proof:

- 1. We pretended w has a fixed sign.
- 2. We integrated over whole space. This is a problem because of boundary terms when integrating by parts.

Now let's be more rigorous. Let $\phi(x) = \frac{x^2}{1+x^2}$, $v = \phi(w) = \phi(f-g)$. Showing w = 0 is equivalent to showing v = 0. We have by the chain rule that

$$\partial_t v - b \partial_h v = 0.$$

Now v has a fixed (positive) sign. This solves item 1.

Denote $L = \|\partial_h f\|_{L^{\infty}} + \|\partial_h g\|_{L^{\infty}} + 1$. Fix $T < \infty$, and study

$$J(t) := \int_{-L(T-t)}^{L(T-t)} v(t,h) dh$$
$$= \int_{-Rt}^{R_t} v$$

Note J is Lipschitz in t, so it has a derivative a.e. By the Fundamental Theorem of Calculus,

$$\partial_t J = \int_{-R_t}^{R_t} \partial_t v - L(v(t, R_t) + v(t, -R_t))$$
$$\int_{-R_t}^{R_t} \partial_t v = \int_{-R_t}^{R_t} b \partial_h v = -\int_{-R_t}^{R_t} (\partial_h b) v + [bv]_{-R_t}^{R_t}.$$

⁴For higher dimensions, the divergence of this vector field has a fixed sign, which says that the flows of the vector field being convergent or divergent.

The term $[bv]_{R_t}^{-R_t}$ may be positive. However, it will be compensated by the – terms because $|b| \leq L$.

$$\partial_t J \le -\int_{-R_t}^{R_t} \underbrace{(\partial_h b)}_{\ge 0} \underbrace{v}_{\ge 0} \le 0$$
$$\partial_t J \le 0.$$

We know J(0)=0 and $J\geq 0$. We conclude that $J\equiv 0$, and v=0 a.e. Because v is Lipschitz, v=0 and f=g.

To be fully rigorous, we should justify interchanging the derivative and integral, and differentiating b (since f, g are only Lipschitz, and may not be twice differentiable).

Exercise 1.3 (for credit): Make this proof rigorous.

Hint: convolve with a function to make it smooth. See the book [Eva10].

What makes the proof work is that the nonlinearity H is convex:

$$\partial_t f - H(\nabla f) = 0.$$

Difficulty comes in when H is not convex or concave.

Any limit point of a solution should be a solution: You can try to run the proof, show the sequence of functions is a Cauchy sequence, and the limit is a solution.

We can write down a formula for the solution.

Proposition 1.4 (Hopf-Lax formula): Let ψ be convex and Lipschitz.⁵ The function

$$f(t,h) = \sup_{h' \in \mathbb{R}} \left(\psi(h - h') - \frac{(h')^2}{4t} \right)$$

is the weak solution of

$$\partial_t f - (\partial_h f)^2 = 0$$
$$f(0, \cdot) = \psi.$$

⁶ Note the last term is related to the convex dual of the nonlinearity: the convex dual of $p\mapsto p^2$ is $q\mapsto \frac{q^2}{4}$. See [Eva10].

Exercise 1.5: Suppose that f solves $\partial_t f - (\partial_h f)^2 = 0$ and $f(0,h) = \psi(h) = \log(e^h + e^{-h})$. Show that for t > 0 small that $\partial_h f(t,0) = 0$. You can use the Hopf-Lax formula. For $t < \infty$ large, $\partial_h^+ f(t,0) > 0 > \partial_h^- f(t,0)$; there is a phase transition in the derivative.

⁵It suffices to be locally semi-convex: for every $\delta > 0$, there exists $C_{\delta} < \infty$ such that for all $\geq \delta$, $h \mapsto f(t,h) + C_{\delta}h^2$. (Note the lower bound is degenerate as $t \to 0$.) More generally, this is true under mild regularity assumptions, though we have to define weak solution differently.

⁶Is there a modification for the pre-limit PDE in F_N ? There would be some Brownian motion. f will be the some expectation of some exponential of Brownian motion. For our application, we will not have a completely closed formula for F_N , so we cannot write a formula of this form before passing to the limit.

1.3 Back to Curie-Weiss

In general it's not true that convergence of functions implies convergence of their derivatives, so we need to prove the following.

Proposition 1.6: If (t,h) is a point of differentiability (in h) of f, and if $F_N \to f$ (pointwise), then

$$\partial_h F_N(t,h) \to \partial_h f(t,h).$$

Proof. F_N is convex in h, so

$$F_N(t, h') \ge F_N(t, h) + \underbrace{\partial_h F_N(t, h)}_{\text{bounded}} (h' - h)$$

Because the derivatives are bounded, we can take a subsequence along which $\partial_h F_N(t,h) \to p$. Then

$$f(t, h') \ge f(t, h) + p(h' - h),$$

and p must be $\partial_h f(t,h)$.

Convergence of F_N . We have

$$\partial_t F_N - (\partial_h F_N)^2 = \frac{1}{N} \partial_h^2 F_N.$$

We want F_N to be close to the HJ solution. Let $w = F_N - f$; then

$$\partial_t w - b\partial_h w = \frac{1}{N} \partial_h^2 F_N$$

where $b = \partial_h F_N + \partial_h f$.

Let $v = \phi(w)$. Before, the difference solved the same equation, but now we have to be more careful.

$$\partial_t v - b\partial_h v = \phi'(w) \frac{1}{N} \partial_h^2 F_N.$$

The argument is similar, but the RHS is different here. Define

$$J(t) = \int_{-R_t}^{R_t} v(t, h) \, dh.$$

We have an extra term

$$\partial_t J \leq \underbrace{\int_{-R_t}^{R_t} b \partial_h v - L(v(t, R_t) + v(t, -R_t))}_{\leq 0} + \underbrace{\int_{-R_t}^{R_t} \underbrace{\phi'(w)}_{\leq 1} \frac{1}{N} \underbrace{\partial_h^2 F_N(t, h)}_{\geq 0} dh}_{}$$

$$\partial_t J \leq \frac{1}{N} \int_{-R_t}^{R_t} \partial_h^2 F_N(t,h) \, dh = \frac{1}{N} [\partial_h F_N]_{-R_t}^{R_t} \leq \frac{2}{N}.$$

where $|\partial_h F_N| \leq 1$ follows from (1). Recall J(0) = 0. So $J(t) \leq \frac{2t}{N}$.

Exercise 1.7: Clean up this proof! How do you get pointwise convergence?

In the rank-1 case, we find some quantities similar to F_N . We have a similar situation with a function which satisfies a similar equation, and want to show it converges to the true solution. We need some L^1 estimate.

Key point: If there are "error terms on the right hand side," we need to estimate it in L^1 in the h variable uniformly in h (locally). We want a "local L_t^{∞} , L_h^1 estimate."

2020/5/21 Lecture 3

1.4 Matrix estimation problem

We saw how to justify the solution in the limit for the Curie-Weiss model, showing error term disappears. Now we consider the matrix estimation problem. This has already been solved by other people. The first that gave a complete solution of this problem is [LM19]. Shortly after, [BM19] gave another proof.

We consider the problem in [Mou18] using the approach in [Mou19].

Problem: Let $\overline{x} = (\overline{x}_1, \dots, \overline{x}_N)$ be a vector of bounded independent random variables with law $P_N = P^{\otimes N}$. For some t > 0, we observe

$$Y = \sqrt{\frac{2t}{N}} \overline{x} \overline{x}^{\top} + W$$

where $W = (W_{ij})$ are independent standard Gaussians. (We don't assume symmetry.) For example, we want to understand

$$\operatorname{mmse}_{N}(t) = \frac{1}{N^{2}} \inf_{\widehat{\theta}} \mathbb{E}\left[|\overline{x}\overline{x}^{\top} - \widehat{\theta}(Y)|^{2}\right]$$
$$= \frac{1}{N^{2}} \mathbb{E}\left[|\overline{x}\overline{x}^{\top} - \mathbb{E}[\overline{x}\overline{x}^{\top}|Y]|^{2}\right].$$

where infimum over all estimators. We assume we can find the best possible estimator given the data, which is the conditional expectation given Y.

Informally,

$$\mathbb{P}\left[\overline{x} = x \text{ and } Y = y\right] = dP_N(x) dy \exp\left(-\frac{1}{2}\left|Y - \sqrt{\frac{2t}{N}}xx^{\top}\right|^2\right)$$

Now compute the conditional law,

$$\mathbb{P}[\overline{x} = x | Y] = \frac{\exp\left(-\frac{1}{2} \left| Y - \sqrt{\frac{2t}{N}} x x^{\top} \right|^{2}\right) dP_{N}(x)}{\int \exp\left(-\frac{1}{2} \left| Y - \sqrt{\frac{2t}{N}} x' x'^{\top} \right|^{2}\right) dP_{N}(x')}$$

Expand the exponent and remove $-\frac{1}{2}|Y|^2$ (which doesn't depend on t,x) to obtain the following.

Definition 1.8: Define

$$H_N^{\circ}(t,x) = \sqrt{\frac{2t}{N}} Y \cdot xx^{\top} - \frac{t}{N} |x|^4.$$

Here Y is the sum of 2 terms, $\sqrt{\frac{2t}{N}}\overline{x}\overline{x}^{\top} + W$. Substituting and expanding gives

$$H_N^{\circ}(t,x) = \sqrt{\frac{2t}{N}} x \cdot W x + \frac{2t}{N} (x \cdot \overline{x})^2 - \frac{t}{N} |x|^4.$$

The first term is the important term; it is like the spin glass model. The other terms will help us. They aren't negligible, but they are here to make out life easier. In summary, the conditional law of $\overline{x}|Y$ has the form of a Gibbs measure, with the most important part looking like the spin glass case.

Denote

$$\langle f(x) \rangle = \frac{\int f(x) \exp(H_N^{\circ}(t, x)) dP_N(x)}{\int \exp(H_N^{\circ}(t, x)) dP_N(x)}.$$

The important difference with the Curie-Weiss model is that $H_N^{\circ}(t,x)$ is still random.

Exercise 1.9 (for credit): Make this rigorous: Show that

$$\langle f(x) \rangle = \mathbb{E}[f(\overline{x})|Y].$$

Why is this model, which looks like the spin glass model, simpler than the spin glass model? The important property is the following.

Proposition 1.10 (Nishimori):

$$\begin{split} \mathbb{E} \left\langle f(x) \right\rangle &= \mathbb{E}[f(\overline{x})] \\ \mathbb{E} \left\langle f(x)g(x') \right\rangle &= \mathbb{E}[\left\langle f(x) \right\rangle \left\langle g(x) \right\rangle] \\ &= \mathbb{E}[\left\langle f(x) \right\rangle \mathbb{E}[g(\overline{x})|Y]] \\ &= \mathbb{E} \left\langle f(x)g(\overline{x}) \right\rangle. \end{split}$$

where x' is an independent copy of x under $\langle \cdot \rangle$. More generally,

$$\mathbb{E} \left\langle f(x, x') \right\rangle = \mathbb{E} \left\langle f(x, \overline{x}) \right\rangle$$

What is x and why is it different from \overline{x} ? Some intuition: Suppose that if we look at Y, we learn nothing about \overline{x} ; then x is a resample. If Y reveals perfect information on \overline{x} , then $x = \overline{x}$. The general case is like an interpolation between these trivial cases.

The independent copy x' is called a **replica**.

This is not trivial, it's not saying $x = \overline{x}$. It is only true because I average outside.

This will help us identify the partial differential equation; we can link it back to the original signal.

We will identify the PDE and compute the mmse by looking at the derivative in the limit, you can recover the phase transition.

Here \overline{x} is the original random vector, and x is our best guess for \overline{x} given what we've seen.

We want to take the log and look for a PDE. But I only have t in this function. If I describe the Curie-Weiss model and had not thought of adding the h; then I would be stuck, I can differentiate in t many times and not find a equation. The relation was between derivatives in t and h. I have t, but what do I relate it to? We play the same game, add a h parameter. Sufficient to relate derivatives.

We have a quadratic term; I would like to add a linear term in x in the model. We want to do this with a constraint: I don't want to destroy the Nishimori property. For Curie-Weiss this was easy, but for spin glasses it's subtle, not easy to figure out in general.

We need to enrich the model somehow!

It has to be rich enough to relate derivatives, and simple enough to compute what happens. Suppose we also observe,

$$\sqrt{2h}\overline{x} + z$$

where z is a standard Gaussian vector. Now I recompute the conditional law of \overline{x} given this rich observation,

$$H_N(t, h, x) = H_N^{\circ}(t, x) + \sqrt{2h}x \cdot z + 2hx \cdot \overline{x} - h|x|^2.$$

The important part is the term $\sqrt{2h}x \cdot z$.

Now I can define the free energy,

$$F_N(t,h) = \frac{1}{N} \mathbb{E} \log \int \exp(H_N(t,h,x)) dP_N(x)$$
$$\overline{F}_N(t,h) = \mathbb{E} F_N(t,h)$$

I take the expectation because the expression inside depends on the randomness in the noise.

Theorem 1.11. $\overline{F}_N \to f$ (local uniform convergence) which is a solution of

$$\partial_t f - (\partial_h f)^2 = 0$$
 on $\mathbb{R}_+ \times \mathbb{R}_+$
 $f(0, \cdot) = F_1(0, \cdot)$

Note $\overline{F}_N(0,\cdot)$ does not depend on N.

1.5 Main steps

The proof has two parts. First we relate the derivates to a variance.

Proposition 1.12:

$$\partial_t \overline{F}_N - (\partial_h \overline{F}_N)^2 = \frac{1}{N^2} \mathbb{E} \left\langle (x \cdot \overline{x} - \mathbb{E} \left\langle x \cdot \overline{x} \right\rangle)^2 \right\rangle.$$

Next, we find a way to express this variance in terms of other things. In the Curie-Weiss model we also described this in terms of derivatives. Here we don't have an equality.

Proposition 1.13: We have

$$\mathbb{E}\left\langle (x \cdot \overline{x} - \mathbb{E}\left\langle x \cdot \overline{x}\right\rangle)^2 \right\rangle \le \frac{1}{N} \partial_h^2 \overline{F}_N + \mathbb{E}[(\partial_h F_N - \partial_h \overline{F}_N)^2]$$

The last term is still random, so we have to control it.

I will justify why this is true, and how to conclude the main result. For the first part, we use the following.

Lemma 1.14.

$$\partial_h \overline{F}_N = \mathbb{E} \left\langle \frac{x \cdot \overline{x}}{N} \right\rangle$$
$$\partial_t \overline{F}_N = \mathbb{E} \left\langle \left(\frac{x \cdot \overline{x}}{N}\right)^2 \right\rangle.$$

The key ingredient is Itô calculus without Itô. Letting X be a standard gaussian,

$$\mathbb{E}[\exp(\sqrt{2t}X - t)] = 1.$$

(Taking the expectation of Brownian motion and subtracting t, we get a constant. Here $\sqrt{2t}X$ is like B_{2t} .) Differentiating with respect to t we should get 0.

$$\mathbb{E}\left[\left(\frac{1}{\sqrt{2t}}X - 1\right)\exp(\sqrt{2t}X - t)\right] = 0.$$

How to see this without Itô calculus? It's something about gaussians. $x \exp(-\frac{x^2}{2})$ want to come together and be integrated out. Using integration by parts,

$$\int x \exp(\sqrt{2t}x - t) \exp\left(-\frac{x^2}{2}\right) dx = \int \sqrt{2t} \exp(\sqrt{2t}x - t) \exp\left(-\frac{x^2}{2}\right) dx.$$

This is Gaussian integration by parts.

Exercise 1.15 (for credit): 1. Prove Lemma 1.14 using Gaussian integration by parts and Nishimori.

2. Relate mmse with $\partial_t \overline{F}_N$.

For Proposition 1.13, letting A be a random variable, notice that

$$\mathbb{E}\left\langle (A - \mathbb{E}\left\langle A\right\rangle)^2\right\rangle = \mathbb{E}\left\langle (A - \left\langle A\right\rangle)^2\right\rangle + \mathbb{E}[(\left\langle A\right\rangle - \mathbb{E}\left\langle A\right\rangle)^2].$$

We do expectation one step at a time. The first term looks like Curie-Weiss, and produces $\frac{1}{N}\partial_h^2\overline{F}_N$. The second term is new, and produces $\partial_h F_N - \partial_h \overline{F}_N$.

The new ingredient is a classical concentration estimate. The Efron-Stein inequality shows $\mathbb{E}[|F_N - \overline{F}_N|^2]$ is small.

Phase transition in t: For small t, we have $\mathbb{E}[\overline{x}_1] = 0$, $\partial_t f = 0$; then it detaches itself from 0, $\partial_t f > 0$. Before the transition, you have maximal error; then the error reduces.

We've talked about best estimation when there are unlimited resources? People believe there is another regime where it is possible but hard to learn something about \overline{x} .

Convergence rate: In the Curie-Weiss, we got the rate for free. In this case, if you follow through, you can get rates. In the paper [Mou19] I was not careful. It's not straightforward to reconstruct the argument in this talk for paper. In the finite-rank case it's not clear that \overline{F}_N is convex. I only know lower bound on convexity, which blows up when $h \to 0$. In the rank-1 case it's convex. From the argument here ,you can get good, N^{-a} convergence (check integration by parts). Certainly you can get $N^{-1/4}$ without a problem.

For p spin, we consider $\partial_t f - (\partial_h f)^p$ instead. Perhaps you can get $|\partial_t f - (\partial_h f)^3| \leq \frac{C}{N} (\partial_h^2 f)^{1/2}$.

2 Critical and near-critical percolation (Gady Kozma)

Critical and near-critical percolation is well-understood in dimension 2 and in high dimensions. The behaviour in intermediate dimensions (in particular 3) is still largely not understood, but in recent years there was some progress in this field, with contributions by van den Berg, Cerf, Duminil-Copin, Tassion and others. We will survey this recent progress (and a few older but not sufficiently known results).

Prerequisites: The Fortuin-Kasteleyn-Ginibre (FKG) and van den Berg-Kesten (BK) inequalities. See e.g. Chapter 2 of Duminil-Copin's lecture notes https://www.ihes.fr/~duminil/publi/2017percolation.pdf or almost any book on percolation theory.

Slides: https://www.math.ubc.ca/Links/OOPS/slides/Kozma_1.pdf

2020/5/25 Lecture 1

Let's start with basic definitions. Examine the graph \mathbb{Z}^d for $d \geq 2$, with edges between nearest neighbors (in L^1 norm). I'm most interested in dimension 3.

For $p \in [0,1]$, keep every edge with probability p and delete it withprobability 1-p, independently for each edge. For background, see Grimmett's book. (The book is a little outdated; I'll discuss results not in the book.)

There is some $p_c \in (0,1)$ (critical p) such that for $p < p_c$, all components ("clusters") of the resulting graph are finite, while for $p > p_c$, there is a unique infinite cluster. Existence of p_c is a consequence of the 0-1 law.

This course is about what happens at p_c . The behaviour at and near p_c is not well understood, except if d = 2 or d > 6. We're interested in the statistics of finite clusters, which has very peculiar properties in all cases we know.

We'll focus on recent advances. The Aizenman-Kesten-Newman argument from the 80's, originally used to prove uniqueness, plays an important role. The argument is in a paper of Cerf in 2015. This is an important piece to understand what is happening at criticality.

What's charming about percolation theory, is that you have to say what is happening at p_c without knowing what the value of p_c is. From a combinatorial point of view: you have some quantity which has an exponential and polynomial correction; you want to understand the polynomial correction without knowing the exponential part. This seems impossible.

Theorem 2.1. $\mathbb{E}_{p_c}(|\mathscr{C}(0)|) = \infty$.

Notation: 0 is the point with all coordinates 0, $\mathcal{C}(x)$ is the cluster containing x. The expected size of the cluster is infinity. Even if it is always finite, it has large tails—it has no finite first moment.

Proof. Fix p and denote $\chi = \mathbb{E}_p(|\mathscr{C}(0)|)$. Let

$$\varepsilon < \frac{1}{4d\chi}$$

We will show that at $p+\varepsilon$ there is no infinite cluster. Consider $p+\varepsilon$ percolation as if we (1) take p-percolation and (2) then "sprinkle" each edge with probability ε . For a vertex x and a sequence of directed edges e_1, \ldots, e_n , denote by E_{x,e_1,\ldots,e_n} the event that 0 is connected to x by a path γ_1 in p-percolation from 0 to e_1 then e_1 is sprinkled, then there is a path γ_2 from e_1 to e_2 then e_2 is sprinkled and soon. We end with a path γ_{n+1} from e_n to x. We require all the γ_i to be disjoint. Clearly $0 \leftrightarrow x$ in $p + \varepsilon$ -percolation if and only if there exist some e_1, \ldots, e_n (possibly empty) such that E_{x,e_1,\ldots,e_n} hold. Then $(0 \leftrightarrow x) = \bigcup_{n=0}^{\infty} \bigcup_{e_1,\ldots,e_n} E_{x,e_1,\ldots,e_n}$ and by the union bound

$$\mathbb{P}_{p+\varepsilon}(0 \leftrightarrow x) \le \sum_{n=0}^{\infty} \sum_{e_1,\dots,e_n} \mathbb{P}(E_{x,e_1,\dots,e_n}).$$

This is the probability in the setting where edge has 3 states: p, ε , or closed (missing). By the BK inequality, this is

$$\leq \sum_{n=0}^{\infty} \sum_{e_1,\dots,e_n} \mathbb{P}_p(0 \leftrightarrow e_1^-) \mathbb{P}_p(e_1^+ \leftrightarrow e_2^-) \cdots \mathbb{P}_p(e_n + \leftrightarrow x) \varepsilon^n.$$

(The paths are disjoint witnesses.)

Summing over all x gives

$$\chi(p+\varepsilon) \leq \sum_{n=0}^{\infty} \varepsilon^n \sum_{e_1,\dots,e_n} \mathbb{P}_p(0 \leftrightarrow e_1^-) \mathbb{P}_p(e_1^+ \leftrightarrow e_2^-) \cdots \mathbb{P}(e_n^+ \leftrightarrow x).$$

(The number of x that e_n^+ is connected to is the size of |(0)|.) Summing over x gives one $\chi(p)$ which we can take out of the sum

$$\sum_{n=0}^{\infty} \varepsilon^n \chi(p) \sum_{e_1, \dots, e_n} \mathbb{P}_p(0 \leftrightarrow e_1^-) \mathbb{P}_p(e_1^+ \leftrightarrow e_2^-) \cdots \mathbb{P}(e_{n-1} + \leftrightarrow x).$$

Even though e_n^+ does not appear in the summand we are summing over it. e_n^+ has 2d possibilities. Summing over e_n^- gives another χ term. Taking both out of the sum gives

$$= \sum_{n=0}^{\infty} \varepsilon^n \cdot 2d\chi(p)^2 \sum_{e_1,\dots,e_{n-1}} \mathbb{P}_p(0 \leftrightarrow e_1^-) \cdots \mathbb{P}_p(e_{n-2}^+ \leftrightarrow e_{n-1}^-)$$
$$= \sum_{n=0}^{\infty} \varepsilon^n (2d)^n \chi(p)^{n+1} < \infty.$$

This shows $p + \varepsilon_c$, contradiction.

(Set of p's for which it is finite is an open set, so at the boundary of the set χ cannot be finite. We used finiteness at the end of the proof.)

The argument also gives

$$\chi(p) \ge \frac{1}{4d(p_c - p)}$$

for all $p < p_c$.

van den Berg-Kesten inequality: Let E be an event on $\{\pm 1\}^n$. We say that an $A \subseteq \{1, \ldots, n\}$ is a witness for E (this is an event itself, denote by ω the parameter) if any ω' such that $\omega'_i = \omega_i$ for all $i \in A$ satisfies that $\omega' \in E$. For example, for the event $0 \to x$, any path between 0 and x can serve as a witness.

For two events E and F we denote by $E \circ F$ the event that they have disjoint witnesses. Then

$$\mathbb{P}(E \circ F) \le \mathbb{P}(E)\mathbb{P}(F).$$

(van den Berg-Kesten 1985 for increasing E and F, Reimer 1997 for arbitrary E and F)

We saw we can argue at p_c using a argument by contradiction. We'll see another example of this argument.

For a set $S \subset \mathbb{Z}^d$ denote by ∂S the set of $x \in S$ with a neighbour $y \notin S$.

Theorem 2.2. Let $S \subset \mathbb{Z}^d$ be some finite set containing 0. Then

$$\sum_{x \in \partial S} \mathbb{P}_{p_c}(0 \stackrel{S}{\longleftrightarrow} x) \ge 1.$$

Proof sketch. Let $x \in \mathbb{Z}^d$. If $0 \leftrightarrow x$ then there exists $0 = y_1, \dots, y_n = x$ such and open paths γ_i such that

- 1. γ_i is from y_i to y_{i+1} and is contained in $y_i + S$.
- 2. The γ_i are disjoint.

We have $n \ge r|x|$ for some number r > 0 that depends on S. (Here |x| is L^1 norm.) A calculation similar to the previous proof shows that

$$\mathbb{P}(0 \leftrightarrow x) \le \sum_{n \ge r|x|} \left(\sum_{y \in \partial S} \mathbb{P}_{p_c}(0 \stackrel{S}{\longleftrightarrow} y) \right)^n.$$

If the value in the parenthesis is smaller than 1, then $\mathbb{P}(0 \leftrightarrow x)$ decays exponentially in |x|, contradicting the previous theorem.

A full proof can be found in H. Duminil-Copin and V. Tassion [DT16].

Theorem 2.3 (Menshikov, Aizenman-Barsky). For any $p < p_c$, $\chi(p) < \infty$.

This is a deeper theorem.

We need to know at the first step that there's one phase transition. Otherwise, there could be p_1, p_2 , such that below p_1 all clusters are finite and expectations are finite, between p_1 and p_2 there is all clusters are finite but expectations are infinite, and p_2 , there is an infinite cluster with positive probability. That would be a completely different picture!

What happens for d > 6? Remote paths of clusters develop independently, as if the other doesn't exist. There's a lot of space. They never get to meet.

2020/5/27 Lecture 2: The Aizenman-Kesten-Newman argument

Today we'll be getting out of the '80's (though this argument is from the 80's).

We start with the idea of using exploration. (This is the one most important takeaway from today's lecture.) We demonstrate this idea with a simple lemma. Recall $\mathscr{C}(0)$ is the cluster containing 0 (everything connected to 0 via open edges).

Lemma 2.4. Let E be the number of open edges in C(0) and let B be the number of closed edges in the boundary. There are C and c such that for any n,

$$\mathbb{P}_p(B+E \le n, |(1-p)E-pB| > \lambda \sqrt{n}) \le Ce^{-c\lambda^2}.$$

The important part is the (1-p)E-pB part. Consider $p=\frac{1}{2}$, so we have roughly the same amount of open and closed edges.

Proof. We define sets of edges $\phi = S_0 \subset S_1 \subset \cdots$ for $i \leq n$ as follows. Assume at step i there exists some edge $e \notin S_i$ such that there is an open path in S_i from 0 to one of the vertices of e. We choose one such e arbitrarily (e.g., by lexicographic order) and define $S_{i+1} := S_i \cup \{e\}$. (Expose the cluster one edge at a time.) If no such e exists (and this happens when $|S_i| = B + E$), let $S_{i+1} = S_i$.

Let

$$X_i = (1 - p)$$
 (#open edges in S_i) – p (#closed edges in S_i).

The key point is X_i is a martingale $(\mathbb{E}[X_{i+1}|X_i]=0)$.

The lemma follows from Azuma-Hoeffding.

Theorem 2.5 (Azuma-Hoeffding). If X_i is a martingale and if $|X_{i+1} - X_i| \le m$ for some numbers m_i , then for any M,

$$\mathbb{P}(|X_n - X_0|) \le 2 \exp\left(\frac{-M^2}{2\sum_{i=1}^N m_i^2}\right).$$

This is a flexible argument. You can start from a set of vertices (not just one), or you can add additional stopping conditions. I'll show one variation.

Lemma 2.6. Let $S \subset \Lambda$ be the set of vertices connected to the boundary. Let E be the number of open edges between vertices of S and let B be the number of closed edges with at least one vertex in S and both vertices in Λ . Let X = (1 - p)E - pB. Then

$$\mathbb{P}(|X| > \lambda n^{d/2}) \le Ce^{-c\lambda^2}.$$

Start from the boundary and examine edges inside the box in arbitrarily order until I discover all vertices connected to the boundary by open paths.

Definition 2.7: Let A, B be subsets of $E \subseteq \mathbb{Z}^d$. We denote by

$$A \stackrel{\stackrel{E}{\longleftrightarrow}}{\longleftrightarrow} B$$

the event that there are 2 disjoint clusters in E which intersect both A and B. We write $A \stackrel{E}{\leftrightarrow} \partial E$.

This is not the same as having 2 disjoint paths. It's stronger. Not only are they disjoint, they are in disjoint clusters.

Theorem 2.8. Let V be the number of edges (x,y) in Λ_n such that $\{x,y\} \stackrel{\leftrightarrow}{\leftrightarrow} \partial \Lambda_n$, i.e., both x and y are connected to $\partial \Lambda_n$ but $x \stackrel{\Lambda_n}{\nleftrightarrow} y$. Then $\mathbb{E}(V) < Cn^{d-\frac{1}{2}}\sqrt{\log n}$.

Clusters can meet outside Λ_n . This proof is a simplification due to Gandolfi-Grimmett-Russo [GGR88] (a nice 4-page paper).

Proof. For $S \subseteq \Lambda_n$ define X(S) to be

(1-p) (# open edges between 2 vertices of S) - p (# closed edges with at least 1 vertex in S, both in Λ_n)

Let C_1, C_2, \ldots be all the clusters in Λ_n that touch the boundary. Then

$$X\left(\bigcup_{i} C_{i}\right) - \sum_{i} X(C_{i}) = pV.$$

This formula is the essence of the proof.

- For an open edge: It's necessarily in one cluster. It contributes 1-p in both terms, which cancel out.
- \bullet For a closed edge with both points in the same cluster: It contributes p to both terms.
- What's left is the event that the closed edge goes between 2 clusters: It is counted twice in the sum, so contributes -p + 2p = p.

The exploration argument shows that with high probability

$$\left| X \left(\bigcup C_i \right) \right| < C n^{d/2} \sqrt{\log n}$$
 $\left| X \left(C_i \right) \right| < C \sqrt{\left| C_i \right|} \sqrt{\log n}$

for all i.

By Cauchy-Schwarz,

$$\sum_{i} \sqrt{|\mathcal{C}_i|} \le \sqrt{\sum_{i} |\mathcal{C}_i|} \sqrt{\sum_{i} 1} \le \sqrt{n^d} \sqrt{n^{d-1}} = n^{d-\frac{1}{2}}.$$

(The number of clusters is bounded by the size of the boundary.) With high probability can be made to mean "with probability $> 1 - n^{-1/2}$ " and we are done.

You can get an estimate for a single edge, by using the theorem for 2n.

Corollary 2.9. For x a neighbour of θ ,

$$\mathbb{P}(\{0, x\} \stackrel{\leftrightarrow}{\leftrightarrow} \partial \Lambda_n) < C\sqrt{\frac{\log n}{n}}.$$

By changing from where you explore you can get all kinds of results. For example, if L is the union of all clusters reaching the left side of Λ_n and R is the union of all clusters reaching the right side of Λ_n then

$$X(L \cup R) - X(L) - X(R)$$

teaches something about edges connected to both the left and the right. Hutchcroft has a version where one explores from random points.

Theorem 2.10. Let $S \subset \mathbb{Z}^d$ be some finite set containing 0. Then $\sum_{x \in \partial S} \mathbb{P}_{p_c}(0 \stackrel{S}{\longleftrightarrow} x) \geq 1$.

Two applications:

Lemma 2.11 (K-Nachmias, 2011). For any $x \in \partial \Lambda_n$, $\Lambda_n := [-n, n]^d$,

$$\mathbb{P}_{p_c}(0 \stackrel{\Lambda_n}{\longleftrightarrow} x) \ge c \exp(-C \log^2 n).$$

Lemma 2.12 (Cerf, 2015). For any $x, y \in \Lambda_n$,

$$\mathbb{P}_{n_c}(x \stackrel{\Lambda_{2n}}{\longleftrightarrow} y) \ge cn^{-C}$$

We can take $C = 2d^2 - 2d$.

All constants c, C might depend on the dimension.

Proof. Assume first that x, y are on the same line (and the distance is even): $x - y = (2k, 0, ..., 0), k \le n$. By the theorem there exists $z \in \partial \Lambda_k$ such that

$$\mathbb{P}(0 \stackrel{\Lambda_k}{\longleftrightarrow} z) \ge \frac{1}{2d|\partial \Lambda_k|} \ge \frac{c}{k^{d-1}}.$$

By rotation and reflection symmetry we may assume z is in some face of Λ_k , for example $z_1 = k$. Let \overline{z} be the reflection of z in the first coordinate i.e. $z = (-z_1, z_2, \dots, z_d)$. By reflection symmetry we also have $\mathbb{P}(0 \stackrel{\Lambda_k}{\longleftrightarrow} \overline{z}) \geq ck^{1-d}$. Translating z to x and \overline{z} to y gives

$$\mathbb{P}(x \overset{x+\Lambda_k}{\longleftrightarrow} x+z), \mathbb{P}(y \overset{y+\Lambda_k}{\longleftrightarrow} y+\overline{z}) \geq \frac{c}{k \cdot d-1}.$$

But $x + z = y + \overline{z}$.

Since $x + \Lambda_k \subset \Lambda_{2n}$ and ditto for $y + \Lambda_k$, we can write

$$\mathbb{P}(x \stackrel{\Lambda_{2n}}{\longleftrightarrow} x + z), \mathbb{P}(y \stackrel{\Lambda_{2n}}{\longleftrightarrow} y + \overline{z}) \ge \frac{c}{k^{d-1}}.$$

By FKG,

$$\mathbb{P}(x \overset{\Lambda_{2n}}{\longleftrightarrow} y) \ge \mathbb{P}(x \overset{\Lambda_{2n}}{\longleftrightarrow} x + z), \mathbb{P}(y \overset{\Lambda_{2n}}{\longleftrightarrow} y + \overline{z}) \ge \frac{c}{k^{2d-2}}.$$

With a slightly smaller c, we can remove the requirement that the distance between x and y is even. If they are not on a line, we define

$$x = x_0, \ldots, x_d = y$$

such that each couple x_i, x_{i+1} differ by only one coordinate. Hence $\mathbb{P}(x_i \overset{\Lambda_{2n}}{\longleftrightarrow} x_{i+1}) \geq cn^{2-2d}$. Using FKG again gives

$$\mathbb{P}(x \stackrel{\Lambda_{2n}}{\longleftrightarrow} y) \ge \mathbb{P}(x_0 \stackrel{\Lambda_{2n}}{\longleftrightarrow} x_1, \cdots, x_{d-1} \stackrel{\Lambda_{2n}}{\longleftrightarrow} x_d)$$
$$\ge \prod_{i=1}^d \mathbb{P}(x_{i-1} \stackrel{\Lambda_{2n}}{\longleftrightarrow} x_i) \ge \frac{c}{n^{2d^2 - 2d}}.$$

Theorem 2.13 (Fortuin-Kasteleyn-Ginibre, Harris). A function $f : \{\pm 1\}^n$ is called increasing if it (weakly) increases in every coordinate. For any 2 increasing functions f, g,

$$\mathbb{E}[fg] \ge \mathbb{E}[f]\mathbb{E}[g].$$

This was recently improved to cn^{-d^2} by van den Berg and Don. Their proof has an interesting topological component (Brouwer's fixed point theorem).

Theorem 2.14 (Cerf, 2015). $\mathbb{P}_{p_c}(\Lambda_{n^c} \stackrel{\leftrightarrow}{\leftrightarrow} \partial \Lambda_n) \leq Cn^{-c}$ for c > 0 small enough. More precisely,

$$\mathbb{P}_{p_c}(\Lambda_{n^{1/(8d^2+8d)-o(1)}} \stackrel{\mbox{\ensuremath{\not\leftarrow}}}{\leftrightarrow} \partial \Lambda_n) \leq C n^{-1/4}$$

Prove using the lemma.

- The theorem actually holds for all p.
- Cerf had a scheme for improving the exponents: repeat the following.
 - 1. Get a better estimate for the number of clusters from $\partial \Lambda_{2n}$ to $\partial \Lambda_n$.
 - 2. Looking at $\sum \sqrt{|\mathcal{C}|}$, get a better estimate for $\mathbb{P}(\{0, x\} \stackrel{\leftrightarrow}{\leftrightarrow} \partial \Lambda)$.
 - 3. Get a better estimate for $\mathbb{P}(\Lambda_{n^c} \stackrel{\leftrightarrow}{\leftrightarrow} \partial \Lambda_n)$. (Use a patching argument.)

Unfortunately, the end result was not a big improvement over $\frac{1}{2}$. (The exponent is $-\frac{2d^2+3d-3}{4d^2+5d-5}$.)

2020/5/28 Lecture 3

3 Branching random walks: some recent results and open questions (Nina Gantert)

We give an introduction to branching random walks and their continuous counterpart, branching Brownian motion. We explain some recent results on the maximum of a branching random walk and its relation to point processes, as well as a connection with fragmentations. The focus will be on open questions.

Preparatory reading: Lyons and Peres, Probability on Trees and Networks http://pages.iu.edu/~rdlyons/prbtree/prbtree.html, Chapters 5.1 (Galton-Watson branching processes) and 13.8 (Tree-indexed random walks)

Further reading:

- Zhan Shi, Branching Random Walks, https://www.lpsm.paris/pageperso/zhan/brw.html
- Julien Berestycki, Topics on Branching Brownian motion, http://www.stats.ox.ac.uk/~berestyc/Articles/EBP18_v2.pdf
- Ofer Zeitouni, Branching Random Walks and Gaussian Fields, http://www.wisdom.weizmann.ac.il/~zeitouni/pdf/notesBRW.pdf

2020/6/1 Lecture 1

There are two ingredients for a branching random walk (BRW):

- 1. Offspring law $p(\cdot)$ with $\sum_{k=1}^{\infty} p(k) = 1$. Assume $\sum_{k=1}^{\infty} kp(k) > 1$. Often we assume p(0) = 0 to simplify.
- 2. Displacement law: Random variable X with Var(X) > 0.

Then

- 1. Start with one particle at 0, particles produce offspring according to $p(\cdot)$.
- 2. Offspring take displacements according to X (all particles behave independently).

How does this crowd of particles behave?

There is a continuous analogue, branching Brownian motion (BBM) (see talk on Friday). It follows Brownian motion and after an exponential lifetime, it splits into 2 particles which continue to follow Brownian motion (and splitting after exponential lifetime, and so on).

We first consider BRW. I will look at it as a tree-indexed random walk in the following sense. Note the Browian motion and branching are independent.

First build a Galton-Watson (GW) process according to $p(\cdot)$. Then label edges with iid random variables distributed as X.

Let D_n be the vertices in generation n. Then $|D_n|$ is the Galton Watson process.

The position of a particle v is the sum of the random variables (displacements) on the edges leading to the vertex v:

$$S_v = \sum_{e \in [0,v]} X_e$$

Recall the following.

Theorem 3.1. If $m := \sum_{k=0}^{\infty} kp(k) > 1$ then $\mathbb{P}(T \text{ infinite}) > 0$.

Here m is the reproduction number.

If p(0) > 0, we can look at $\mathbb{P}^*[\cdot] = \mathbb{P}[\cdot||D_n| > 0, \forall n]$. This event has positive probability. The collection $(S_v)_{v \in D_n}$ are random variables which are not independent.

There is a more general model: Particles produce "offspring and displacements" at once according to some point process.

For example, the point process could be:

- produce particles at position 1 and -1,
- produce 3 particles at positions 3,

each with probability $\frac{1}{2}$. There is dependence between siblings, and displacements are not independent of tree.

Many things I will say will generalize to the general model.

The first question I will address is the behavior of the rightmost particle:

$$M_n = \sup_{v \in D_n} S_v.$$

This is the maximum of variables with are not independent. Assume the exponential moment condition

$$\mathbb{E}[e^{\lambda X}] < \infty$$

for some $\lambda > 0$ and define a large deviation rate function

$$I(y) = \sup_{\lambda} [\lambda y - \log \mathbb{E}[e^{\lambda X}]]$$

It is a fact that for $y > \mathbb{E}[X]$,

$$\frac{1}{n}\log \mathbb{P}[S_n \ge ny] \to -I(y).$$

You can proved one direction guite easily:

$$\mathbb{P}(S_n > ny) \le \mathbb{E}[e^{\lambda S_n}]e^{-\lambda ny}$$
$$= e^{-nI(y)}$$

by Chebyshev's inequality and optimizing the RHS over λ . Define

$$x^* = \sup \left\{ s \ge \mathbb{E}[X] : I(s) \le \log m \right\}.$$

We have the following old result.

Theorem 3.2 (Biggins, Hammersley, Kingman).

$$\frac{M_n}{n} \to x^*$$
 \mathbb{P}^* -a.s.

The method of proof is useful.

Exercise 3.3: Assume the theorem.

- 1. Let $X \stackrel{d}{=} N(0,1)$. Compute x^* .
- 2. Suppose $p(3)=1, \mathbb{P}[X=0]=\frac{1}{2}=\mathbb{P}[X=1].$ Compute x^* . Do we have $\mathbb{P}[M_n=n,\forall n]>0$?
- 3. Same as (ii) if p(2) = 1.

I'll give my favorite proof of the theorem.

Intuition: At time n, we have $\approx m^n$ particles. For each $v \in D_n$, $\mathbb{P}(S_u \ge ny) \approx e^{-nI(y)}$. We have 2 competing effects: the probability decays exponentially but the number of particles increase exponentially. These effects shold balance: if

$$e^{-nI(y)}\underbrace{e^{n\log m}}_{m^n} = 1$$

then $y = x^*$.

Proof. 1. First moment method: This will give an upper bound.

$$\mathbb{P}[M_n \ge ny] \le \mathbb{E}\left[\sum_{v \in D_n} I_{S_v \ge ny}\right]$$
$$= \underbrace{\mathbb{E}[|\Delta_n|]}_{m^n} \underbrace{\mathbb{P}[S_n \ge ny]}_{\le e^{-nI(y)}}.$$

using the independence assumption. Hence, if $I(y) > \log m$, then

$$\sum_{u} \mathbb{P}[M_u \ge ny] < \infty$$

$$\implies \limsup_{n \to \infty} \frac{M_n}{n} \le y.$$

2. Embedded tree: Assume $y < x^*$. Choose $\varepsilon > 0$ such that $I(y) - 2\varepsilon < \log m$. Then using the lower bound

$$\mathbb{P}[S_k \ge ky] \ge e^{-k(I(y) - \varepsilon)}$$

for $k \geq k_0(\varepsilon)$. I don't have strict inequality $e^{-kI(y)}$, I have an ε .

Now I'm doing a kind of percolation:

- keep $v \in D_k$ if $\frac{S_v}{k} \ge y$
- \bullet delete v otherwise
- go at level 2k, etc.

Continue only vertices which I kept at the first step, where the mean is large enough. Now I have an embedded GW-tree $\widetilde{T} \subset T$. If \widetilde{T} is infinite, the maximum is at least y. Question: what is \widetilde{m} ? We have

$$\widetilde{m} = m^k \mathbb{P}[S_k \ge ky]$$

$$> e^{k \log m} e^{-k(I(y) - \varepsilon)} > e^{\varepsilon k} > 1$$

for k large enough. Then

$$\mathbb{P}[\lim_{n\to\infty}\inf\frac{M_n}{n}\geq y]>0.$$

Now we need a 0-1-law for inherited properties. Call a property A of trees **inherited** if each finite trees has A, and if T has A, then all descendant trees of the children of the root have it. Then $\mathbb{P}^*[T \text{ has } A] \in \{0,1\}$. Thus we conclude the probability is 1.

Proof of 0-1 law. We have

$$\begin{split} \mathbb{P}[T \text{ has } A] &= \mathbb{E}[\mathbb{P}[T \text{ has } A||D_1|]] \\ &\leq \mathbb{E}[\mathbb{P}[T^{(1)} \text{ has } A, \dots, T^{(|D_1|)} \text{ has } A]||D_1|] & \text{inherited property} \\ &= \mathbb{E}[\mathbb{P}[T \text{ has } A]^{|D_1|}] & \text{independence} \end{split}$$

Hence

$$\mathbb{P}[T \text{ has } A] \leq f(\mathbb{P}[T \text{ has } A])$$

where $f(s) = \sum_{k=0}^{\infty} s^k p(k) = \mathbb{E}[s^{|D_1|}]$. On the other hand

$$\mathbb{P}[T \text{ has } A] \ge 0$$

$$q := \mathbb{P}[\lim_{n} |D_n| = 0]$$

This means that

$$\mathbb{P}[T \text{ has } A] \in \{q, 1\}.$$

But if we condition, we get

$$\mathbb{P}^*[T \text{ has } A] \in \{0, 1\}$$

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