



**Weierstrass Institute for
Applied Analysis and Stochastics**



A large-deviations principle for all the components in a sparse inhomogeneous Erdős-Rényi graph

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based on joint work (in progress) with Luisa Andreis (Florence), Heide Langhammer and Robert Patterson (WIAS)

The **MARCUS-LUSHNIKOV model** is a non-spatial (i.e., a mean-field) coagulation model [MARCUS 1968], [GILLESPIE 1972], [LUSHNIKOV 1978]:

Continuous-time Markov process of vectors of particle masses at time $t \in [0, \infty)$:

$$M_1^{(N)}(t) \geq M_2^{(N)}(t) \geq M_3^{(N)}(t) \geq \dots \geq M_{n(t)}^{(N)}(t) \geq 1, \quad \sum_{i=1}^{n(t)} M_i^{(N)}(t) = N.$$

We start with $M_i^{(N)}(0) = 1$ for any i .

Coagulation mechanism:

Particles with masses m and \tilde{m} coagulate after an exponential random time with parameter $K_N(m, \tilde{m})$ (the **coagulation kernel**) independently of all the other particles.

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Our main question:

Is there a gelation phase transition at some fixed time $t_c \in (0, \infty)$ in the limit $N \rightarrow \infty$?

That is, is there a (deterministic) time after which a **gel** emerges, i.e., a particle with size $M_1^{(N)}(t) \asymp N$?

Here, we make the special choice of the **multiplicative kernel**:

$$K_N(m, \tilde{m}) = \frac{m \tilde{m}}{N}.$$

Advantage:

The model is now a function of a time-dependent version of the well-known **ERDŐS-RÉNYI random graph** model. Indeed, the vector $(M_i^{(N)}(t))_{i=1}^{n(t)}$ is in distribution equal to the collection of sizes of all the connected components of the graph $\mathcal{G}(N, 1 - e^{-t/N})$.

Explanation:

Equip each unordered pair $\{i, j\}$ of different numbers in $\{1, \dots, N\}$ independently with an exponentially distributed random time $e_{i,j}$ with expected value N . After the elapsure of $e_{i,j}$, there is a bond created between i and j . Then, at time t , for each pair, the probability to have a bond between them is equal to $1 - e^{-t/N}$.

Imagine the component containing i (with size m) and the one containing j (with size \tilde{m}) have been turned into a new component (with size $m + \tilde{m}$). Then the new component inherits all the active $m\tilde{m}$ exponential random times of the two earlier components.

From now, we stick to the sparse Erdős–Rényi graph on $[N] = \{1, \dots, N\}$.

Goal 1: Explicit joint **large-deviation principle** for the statistics of all the component sizes k , distinguished into microscopic ($k \asymp 1$), mesoscopic ($1 \ll k \ll N$) and the macroscopic ($k \asymp N$) sizes. Explicit identification of the gelation phase transition as a consequence.

Goal 2: The same for a "spatial" version, the *inhomogeneous Erdős-Rényi graph*.

Earlier works on LDPs for sparse random graphs:

- [O'CONNELL 1998]: LDP for size of largest component and number of isolated points
- [ENGEL, MONASSON, HARTMANN 2004]: LDP for free energy of a tilted version with weights on the number of components, connections with Potts model.
- [BORDENAVE, CAPUTO 2015]: LDP for the microscopic connected subgraphs
- [PUHALSKII 2005]: LDP for the number of components, number of macroscopic components, number of excess edges in them. (Proof ansatz and rate function very different from ours).

Fix $t > 0$ and consider the standard Erdős–Rényi graph $\mathcal{G}(N, \frac{t}{N})$ with components of sizes $S_1^{(N)} \geq S_2^{(N)} \geq \dots \geq S_n^{(N)} \geq 1$.

Microscopic and macroscopic empirical measures of the particle sizes:

$$\text{Mi}^{(N)} = \frac{1}{N} \sum_{i=1}^n \delta_{S_i^{(N)}} \quad \text{and} \quad \text{Ma}^{(N)} = \sum_{i=1}^n \delta_{\frac{1}{N} S_i^{(N)}}.$$

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Then $\text{Mi}^{(N)}$ is a random member of the set $\mathcal{N} = \mathcal{N}(1)$, where

$$\mathcal{N}(c) = \left\{ \lambda \in [0, \infty)^{\mathbb{N}} : \sum_{k \in \mathbb{N}} k \lambda_k = c \right\} \quad (\text{coordinatewise top.}).$$

$\text{Ma}^{(N)}$ is a random element of $\mathcal{M}_{\mathbb{N}_0} = \mathcal{M}_{\mathbb{N}_0}(1)$, where

$$\mathcal{M}_{\mathbb{N}_0}(c) = \left\{ \alpha \in \mathcal{M}_{\mathbb{N}_0}((0, 1]) : \int_{(0,1]} x \alpha(dx) = c \right\} \quad (\text{vague top.}).$$

and $\mathcal{M}_{\mathbb{N}_0}((0, 1])$ is the set of all measures on $(0, 1]$ with values in \mathbb{N}_0 .

Note that the **total masses**

$$c_\lambda = \sum_{k \in \mathbb{N}} k \lambda_k \quad \text{and} \quad c_\alpha = \int_{(0,1]} x \alpha(dx)$$

are discontinuous functions of λ resp. α .

LDP for the micro- and macroscopic parts

As $N \rightarrow \infty$, the pair $(\text{Mi}^{(N)}, \text{Ma}^{(N)})$ satisfies an LDP with rate function

$$I(\lambda, \alpha; t) = \begin{cases} I_{\text{Mi}}(\lambda; t) + I_{\text{Ma}}(\alpha; t) + (1 - c_\lambda - c_\alpha) \left(\frac{t}{2} - \log t \right), & \text{if } c_\lambda + c_\alpha \leq 1, \\ \infty & \text{otherwise,} \end{cases}$$

where

$$I_{\text{Mi}}(\lambda; t) = \sum_{k=1}^{\infty} \lambda_k \log \frac{k! t \lambda_k}{e k^{k-2}} + c_\lambda \left(1 + \frac{t}{2} - \log t \right),$$

$$I_{\text{Ma}}(\alpha; t) = \int_0^1 \left[x \log \frac{x}{1 - e^{-tx}} + \frac{t}{2} x(1 - x) \right] \alpha(dx).$$

Microscopic total mass phase transition

1. For $t \in (0, 1)$, the minimum of micro-part of the rate function is attained precisely at

$$\lambda_k^*(c; t) = \frac{k^{k-2} c^k t^{k-1} e^{-ctk}}{k!}, \quad k \in \mathbb{N},$$

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2. For $t \in (1, \infty)$, the minimum of the micro total mass rate function is attained precisely at $c = \beta_t$ for some $\beta_t \in (0, \frac{1}{t})$, given as the smallest positive solution to $\log \beta_t = t\beta_t - t$. The infimum is attained precisely at $(\lambda, \alpha) = (\lambda^*(\beta_t; t), (1 - \beta_t, 0, \dots))$.

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Hence, $t_c = 1$ is the gelation transition time. On a linear level, we can say:

- Before time 1, all particles are finitely large, and the statistics of their sizes follow the **Borel distribution**.
- After time 1, there is precisely one macroscopic particle of size $\sim (1 - \beta_t)N$, and a Borel-distributed statistics of remaining particle sizes.

LDP for mesoscopic total mass

Fix $t \in [0, \infty)$ and $\varepsilon > 0$ and $R \in \mathbb{N}$. Then the mesoscopic (ε, R) -total mass,

$$\overline{\text{Me}}_{R,\varepsilon}^{(N)} = \frac{1}{N} \sum_{i: R < M_i^{(N)}(t) < \varepsilon N} S_i^{(N)}.$$

satisfies an LDP with some rate function $\mathcal{J}_{\text{Me}}^{(\varepsilon, R)}$ whose limit for $\varepsilon \downarrow 0$ and $R \rightarrow \infty$ is equal to

$$\mathcal{J}_{\text{Me}}(c) = (1 - c) \left(\log(1 - c)t - \frac{(1 - c)t}{2} \right) + \frac{t}{2} - \log t.$$

- \mathcal{J}_{Me} is strictly increasing and has a unique zero at $c = 0$.
- We also proved that $\overline{\text{Me}}_{R_N, \varepsilon_N}^{(N)}$ satisfies an LDP with rate function \mathcal{J}_{Me} if $1 \ll R_N \leq N\varepsilon_N \ll N$.

Let $\mathbb{P}_{k,p}$ be the probability measure for $\mathcal{G} \sim \mathcal{G}(k, p)$. Put

$$\mu_k(p) = \mathbb{P}_{k,p}(\mathcal{G} \text{ is connected}),$$

then we have

Distribution of statistics

For any N and any $\ell = (\ell_k)_k \in \mathbb{N}_0^{\mathbb{N}}$ satisfying $\sum_k k\ell_k = N$, write

$$A_N(\ell) = \bigcap_{k \in \mathbb{N}} \{\#\{i: S_i^{(N)} = k\} = \ell_k\},$$

then

$$\mathbb{P}_{N,p}(A_N(\ell)) = N! \prod_k \frac{\mu_k(p)^{\ell_k} (1-p)^{\frac{1}{2}k(N-k)\ell_k}}{k!^{\ell_k} \ell_k!}.$$

Proof: elementary combinatorics.

Micro and macro asymptotics [STEPANOV 1970]



$$(1-p)^{\frac{1}{2}(k-1)(k-2)} \leq \frac{\mu_k(p)}{k^{k-2}p^{k-1}} \leq 1, \quad k \in \mathbb{N}.$$

In particular, if $k = o(\sqrt{N})$, then

$$\mu_k\left(\frac{t}{N}\right) = k^{k-2}\left(\frac{t}{N}\right)^{k-1}, \quad N \rightarrow \infty.$$



$$\mu_{\lfloor \alpha N \rfloor}\left(\frac{t}{N}\right) \sim \left(1 - \frac{\alpha t}{e^{\alpha t} - 1}\right) (1 - e^{-\alpha t})^{\alpha N}, \quad \alpha \in (0, 1).$$

Consequences for the coagulation model:

- In [SMOLUCHOWSKI 1916] a system of ODEs is introduced for the evolution of the (microscopic) particle sizes:

$$\frac{d}{dt} \lambda_k(t) = \frac{1}{2} \sum_{m, \tilde{m} \in \mathbb{N}: m + \tilde{m} = k} \lambda_m(t) \lambda_{\tilde{m}}(t) K(m, \tilde{m}) - \lambda_k(t) \sum_{m \in \mathbb{N}} \lambda_m(t) K(k, m),$$

where $K = \lim_{N \rightarrow \infty} N K_N$, and

$$\lambda_m(t) = \lim_{N \rightarrow \infty} \frac{1}{N} \#\{\text{particles at time } t \text{ of size } k\}.$$

- One can check that the minimizers λ_k^* of our variational formula satisfy them.
- Convergence of stochastic coagulation processes towards these ODEs was expected for long time, but the first rigorous proof was given only in [LANG, NGUYEN 1980].
- In [LUSHNIKOV 1978] the formation of a gel is realized and explained.
- Pathwise large deviations appear cumbersome, but doable.
- Such LDPs have been derived by [MIELKE et. al. (2017)] for general chemical reactions, following a Freidlin-Wentzell approach, but the rate function is rather inexplicit and not easy to evaluate at a fixed time.

Consider the non-interacting **Bose gas** in the thermodynamic limit at **temperature** $1/\beta \in (0, \infty)$ with **particle density** $\rho \in (0, \infty)$. Then the partition function is given by

$$Z_N(\beta, \rho) = \sum_{(\ell_k)_{k \in \mathbb{N}} \in \mathbb{N}_0^{\mathbb{N}}: \sum_k k \ell_k = N} \prod_k \frac{N^{\ell_k}}{\ell_k! k^{\ell_k}} [\rho(4\pi\beta k)^{\frac{d}{2}}]^{-\ell_k}.$$

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The free energy per particle is then

$$f(\beta, \rho) = \lim_{N \rightarrow \infty} \frac{1}{N} \log Z_N(\beta, \rho) = - \inf_{\lambda \in \mathcal{N}(\rho)} I(\lambda), \quad \text{where} \quad I(\lambda) = \sum_{k \in \mathbb{N}} \lambda_k \log \frac{\lambda_k k}{(4\pi\beta k)^{\frac{d}{2}} e}.$$

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Comparison: In Lushnikov's model, we face roughly

$$t^N e^{\frac{t}{2}N} \sum_{(\ell_k)_{k \in \mathbb{N}}: \sum_k k \ell_k = N} \prod_k \frac{k^{(k-2)\ell_k} t^{-\ell_k}}{\ell_k! k!^{\ell_k}}.$$

The two respective minimizers are

$$k\lambda_k^{(\text{Lush})}(c; t) = \frac{1}{t} \frac{(cte^{-ct})^k}{k^{1-k} k!} \quad \text{and} \quad k\lambda_k^{(\text{BEC})}(\alpha; t) = \frac{1}{\rho(4\pi\beta)^{\frac{d}{2}}} \frac{e^{-\alpha k}}{k^{\frac{d}{2}}},$$

where c and α control the value of $\sum_k k\lambda_k$ (note that $k^{1-k} k! \asymp k^{3/2}$).

One difference: In the non-interacting Bose gas, the macroscopic part gives no energetic contribution, while in the Lushnikov model it does.

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In the Bose gas, increasing ρ drives more and more particles into the finite cycles. There is a natural threshold, the critical inverse temperature β_c , characterised by

$$(4\pi\beta)^{\frac{d}{2}}\rho = \sum_{k \in \mathbb{N}} k^{-\frac{d}{2}}.$$

Only when all finite cycles are filled entirely, the first “infinite” cycle arises.

The BEC is a **saturation** transition.

In contrast, in the Lushnikov model, increasing t makes each particle larger, until some decide to make the jump to infinity. However, the other micro particles keep growing (recall that $\beta_t < \frac{1}{t}$).

The gelation phase transition is an **explosion** transition.

- **type space:** compact metric space \mathcal{S}
- **vertex distribution:** probability measure μ on \mathcal{S}
- **connectivity probability function:** positive symmetric irreducible kernel κ from \mathcal{S} to \mathcal{S} .
- **vertex set:** $[N] = \{1, \dots, N\}$. Vertex i has the type $x_i \in \mathcal{S}$. Type vector $x = (x_1, \dots, x_N) \in \mathcal{S}^N$.

$\mathcal{G}_N = \mathcal{G}([N], x, \frac{1}{N}\kappa)$ is the graph on $[N] = \{1, \dots, N\}$, having a **bond** $\{i, j\}$ with **probability** $\frac{1}{N}\kappa(x_i, x_j) \wedge 1$, independently over all pairs (i, j) with $i \neq j$.

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- There is a spatial coagulation model that can be mapped onto this graph model.
- The 120-pages article [BOLLOBAS, JANSON, RIORDAN 2007] derived a sufficient and necessary criterion for the **phase transition of the existence of a giant component** in \mathcal{G}_N . The main tool is a **multitype branching process**.
- We will prove the extension to an **LDP for the statistics of the microscopic and the macroscopic components** of \mathcal{G}_N , and obtain this criterion independently in a different way. The main tool is the explicit identification of the joint distribution of the statistics of all the connected components according to their multi-types.

For a measure ν on \mathcal{S} , introduce the operator

$$T_{\kappa, \nu}: L^2(\nu) \rightarrow L^2(\nu), \quad T_{\kappa, \nu} f(x) = \int_{\mathcal{S}} \kappa(x, y) f(y) \nu(dy),$$

and its norm

$$\Sigma(\kappa, \nu) = \|T_{\kappa, \nu}\|_{L^2(\nu)} = \sup_{f \in L^2(\nu): \|f\|_{L^2(\nu)}=1} \|T_{\kappa, \nu} f\|_{L^2(\nu)}.$$

Existence of a giant component

- If $\Sigma(\kappa, \mu) \leq 1$, then the largest component of \mathcal{G}_N has size $o(N)$ as $N \rightarrow \infty$ with high probability (in fact, $O(\log N)$).
- If $\Sigma(\kappa, \mu) > 1$, then it has size $\asymp N$. More precisely, if $\rho: \mathcal{S} \rightarrow [0, \infty)$ denotes the maximal solution of

$$\rho = 1 - e^{-T_{\kappa, \mu} \rho},$$

then the size of the largest component of \mathcal{G}_N is $\sim N \int_{\mathcal{S}} \rho(x) \mu(dx)$.

The sizes of the microscopic clusters are characterized in terms of the distribution of the sizes of the offspring of the **multitype branching process**, in which each particle of type $x \in \mathcal{S}$ has offspring with distribution that is a Poisson process with intensity $\kappa(x, y) \mu(dy)$.

Assume that \mathcal{S} is a **finite set**. Assume that $\mu_N = \frac{1}{N} \sum_{i=1}^N \delta_{x_i} \implies \mu$ as $N \rightarrow \infty$.

We denote by $(\mathcal{C}_i)_i$ the collection of the **connected components** of \mathcal{G}_N .

Let $\eta_r(A)$ be the **number of type- r sites** in $A \subset [N]$, and $\eta(A) = (\eta_r(A))_{r \in \mathcal{S}}$.

For $k \in \mathbb{N}_0^{\mathcal{S}}$ let $\text{Mi}_N(k) := \frac{1}{N} \sum_i \delta_{\eta(\mathcal{C}_i)}(k)$, then $\sum_k \text{Mi}_N(k) k_r = \mu_N(r)$.

For $\alpha \in (0, 1]^{\mathcal{S}}$ let $\text{Ma}_N(k) := \sum_i \delta_{\frac{1}{N} \eta(\mathcal{C}_i)}(k)$, then $\int_{(0,1]^{\mathcal{S}}} \text{Ma}_N(dy) y_r = \mu_N(r)$.

Joint distribution of the cluster types

For any $l = (l_k)_{k \in \mathbb{N}_0^{\mathcal{S}}} \in \mathbb{N}_0^{(\mathbb{N}_0^{\mathcal{S}})}$ satisfying $\sum_k l_k k_r = N \mu_N(r)$ for any $r \in \mathcal{S}$,

$$\mathbb{P}(N \text{Mi}_N(k) = l_k \forall k) = \left(\prod_{r \in \mathcal{S}} (N \mu_N(r))! \right) \\ \times \prod_{k \in \mathbb{N}_0^{\mathcal{S}}} \frac{p_N(k)^{l_k}}{l_k! (k_r!)^{l_k}} \left(\prod_{r,s \in \mathcal{S}} \left(1 - \frac{\kappa(r,s)}{N} \right)^{k_s (N \mu_N(r) - k_r) / 2} \right)^{l_k}$$

with $p_N(k)$ the connection probability for the graph $\mathcal{G}(|k|, x, \frac{1}{N} \kappa)$ for any k -compatible x .

Define

$$\tau(k) := \sum_{T \in \mathcal{T}(k, x)} \prod_{\{i, j\} \in E(T)} \kappa(x_i, x_j), \quad k \in \mathbb{N}_0^{\mathcal{S}},$$

where $x \in \mathcal{S}^{|k|}$ is k -compatible, and $\mathcal{T}(k, x)$ is the set of spanning trees on $[|k|]$.

Notable extension of [STEPANOV 1970]:

Asymptotics of $p_N(k)$ as $N \rightarrow \infty$

$$p_N(k) \sim N^{1-|k|} \tau(k), \quad k \in \mathbb{N}_0^{\mathcal{S}}.$$

and

$$\frac{1}{N} \log p_N(\lfloor Ny \rfloor) \rightarrow \sum_{r \in \mathcal{S}} y_r \log \left(1 - e^{-\sum_{s \in \mathcal{S}} \kappa(r, s) y_s} \right), \quad y \in (0, 1]^{\mathcal{S}}.$$

The second assertion is of independent interest and is also proved for \mathcal{S} a compact metric space. The technical problem is that giant clusters can be connected with just one bond, whose probability is not seen on the exponential scale.

Denote $c_r(\lambda) = \sum_{k \in \mathbb{N}^{\mathcal{S}}} \lambda_k k_r$ and $c_r(\alpha) = \int_{(0,1]^{\mathcal{S}}} \alpha(dy) y_r$.

The LDP

As $N \rightarrow \infty$, the pair (M_{iN}, M_{aN}) satisfies a large-deviations principle with rate function

$$I(\lambda, \alpha) = I_{\text{Mi}}(\lambda) + I_{\text{Ma}}(\alpha) + I_{\text{Me}}(\mu - c(\lambda) - c(\alpha)),$$

where

$$I_{\text{Mi}}(\lambda) = \sum_{k \in \mathbb{N}^{\mathcal{S}}} \lambda_k \log \frac{\lambda_k}{\tau(k) \prod_{s \in \mathcal{S}} \frac{\mu_s^{k_s}}{k_s!}} + \sum_{k \in \mathbb{N}^{\mathcal{S}}} \lambda_k (|k| - 1) + \frac{1}{2} \langle c(\lambda), \kappa \mu \rangle,$$

$$I_{\text{Ma}}(\alpha) = \int_{[0,1]^{\mathcal{S}}} \alpha(dy) \left(\left\langle y, \log \frac{y}{1 - e^{-\kappa * y}} \right\rangle + \frac{1}{2} \langle y, \kappa * (\mu - y) \rangle \right),$$

$$I_{\text{Me}}(\nu) = \left\langle \nu, \log \frac{\nu}{(\kappa * \nu) \mu} \right\rangle + \frac{1}{2} \langle \nu, \kappa * \mu \rangle.$$

- entropies \iff combinatorics
- terms with $\frac{1}{2}$ \iff non-connection probabilities
- term τ times Poisson \iff reference process, conditioned on being connected

- We indeed prove this also for \mathcal{S} a compact metric space. The lift from discrete \mathcal{S} to continuous \mathcal{S} is a cumbersome and technical work in the spirit of the DAWSON–GÄRTNER theorem.
- We do not know about earlier work in that direction.
- One application is to i.i.d. random $x_1, \dots, x_N \implies$ quenched LDP. Annealed versions follow easily.
- Standard consequences are contracted separate LDPs for Mi_N and Ma_N . (\implies interesting conditional phase transition, see later)
- We abstained from formulating an LDP for the mesoscopic part.

Given $c = (c_r)_{r \in \mathbb{N}_0^{\mathcal{S}}}$, the Euler-Lagrange equations for minimizers λ of I_{Mi} subject to $c(\lambda) = c$, i.e., $\sum_k \lambda_k k_r = c_r$ for $r \in \mathcal{S}$, are

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Existence of $\lambda^*(c)$

- $t(c)$ is a solution $\iff \tilde{c} \mapsto t(\tilde{c})$ is invertible and the inverse map is analytic in $t(c)$,
- this invertibility is true if and only if $\Sigma(\kappa, c) < 1$.

Given $c = (c_r)_{r \in \mathbb{N}_0^S}$, the Euler-Lagrange equations for minimizers λ of I_{Mi} subject to $c(\lambda) = c$, i.e., $\sum_k \lambda_k k_r = c_r$ for $r \in \mathcal{S}$, are

$$\lambda_k = \tau(k) \prod_{r \in \mathbb{N}_0^S} \frac{t_r^{k_r}}{k_r!}, \quad k \in \mathbb{N}_0^S.$$

The only candidate is $t_r(c) = c_r e^{-\kappa * c(r)}$. Call the **solution $\lambda^*(c)$** if it exists.

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Minimizers of LDP rate function

- If $\Sigma(\kappa, \mu) \leq 1$, then $\lambda^*(\mu)$ exists, and $(\lambda^*(\mu), 0)$ is the minimizer of I . No giant component arises.
- If $\Sigma(\kappa, \mu) > 1$, then the optimal microcluster distribution c^* is characterised by $c_r = \mu_r e^{\kappa * (\mu - c)(r)}$, and the minimizer of I is equal to $(\lambda(c^*), \delta_{\mu - c^*})$. The latter corresponds to a giant cluster with $\sim N(\mu_r - c_r^*)$ vertices of multitype $r \in \mathbb{N}_0^S$ for any r .

Recall the **multitype branching process**, in which each particle of type $x \in \mathcal{S}$ has offspring with distribution that is a Poisson process with intensity $\kappa(x, y) \mu(dy)$.

Denote by $\Xi(dr)$ the **entire progeny** (total offspring) of type $r \in \mathcal{S}$ the process. Let P_r denote the probability measure when the process starts with just one particle of type r .

Then

$$\mu(dr)P_r(\Xi \in dk) = \lambda_\mu(dk)k(dr), \quad k \in \mathcal{M}_{\mathbb{N}_0}(\mathcal{S}), r \in \mathcal{S}.$$

In words: the empirical statistics of the microscopic components in \mathcal{G}_N in the subcritical case approximate the distribution of the total offspring of the characteristic branching process.

Contraction principle \implies Ma_N satisfies an LDP with rate function

$$\mathcal{I}_{\text{Ma}}(\alpha) = \inf_{\lambda} I(\lambda, \alpha) = I_{\text{Ma}}(\alpha) + J(\mu - c_{\alpha}),$$

where

$$J(c) = \begin{cases} I_{\text{Mi}}(\lambda_c) & \text{if } \Sigma(\kappa, c) \leq 1, \\ I_{\text{Mi}}(\lambda_{b^*}) + I_{\text{Me}}(c - b^*) & \text{if } \Sigma(\kappa, c) > 1, \end{cases}$$

and $b^* = b^*(c) \in \mathcal{M}(\mathcal{S})$ is the minimal solution $\neq c$ of the characteristic equation

$$\kappa(c - b^*)(r) b^*(dr) = (c - b^*)(dr), \quad b^* \leq c,$$

and b^* is saturated in the sense that $\Sigma(\kappa, b^*) = 1$.

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Hence, conditional on $\{\text{Ma}_N \approx \alpha\}$, we have, as $N \rightarrow \infty$,

$$\text{Mi}_N \xrightarrow{N \rightarrow \infty} \begin{cases} \lambda_{\mu - c_{\alpha}} & \text{if } \Sigma(\kappa, \mu - c_{\alpha}) < 1 \implies \text{no mesoscopic part,} \\ \lambda_{b^*} & \text{if } \Sigma(\kappa, \mu - c_{\alpha}) \geq 1 \implies \text{mesoscopic part.} \end{cases}$$

\implies **saturation phase transition**: If the macroscopic part α is fixed, and more and more bonds are thrown in, then first the microscopic part increases until λ_{b^*} is attained, then it is frozen, and only the mesoscopic part increases. (\implies **frozen percolation**)