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Statistical Physics of Complex Networks

From Graph Ensembles to Emergent Phenomena

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Statistical Physics of Complex Networks

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*To Marco and Sofia,
who taught me that every connection matters.*

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Preface

THIS monograph grew out of a series of graduate lectures I delivered at the Institute for Advanced Study between 2019 and 2023. The central premise is deceptively simple: many of the most challenging problems in modern physics — from the glass transition to quantum decoherence — can be reframed as questions about the structure and dynamics of complex networks.

The approach taken here is deliberately interdisciplinary. Readers will encounter ideas drawn from statistical mechanics, graph theory, information geometry, and the theory of disordered systems. Each chapter begins with the necessary background and builds toward results at the research frontier. I have aimed for a self-contained treatment that serves both as a graduate-level textbook and as a reference for active researchers.

Princeton, New Jersey

January 2025

Part I

Foundations

The most exciting phrase to hear in science, the one that heralds new discoveries, is not “Eureka!” but “That’s funny . . .”

Isaac Asimov

Chapter 1

The Landscape of Network Science

THE study of networks has undergone a remarkable transformation over the past two decades. What began as a modest offshoot of graph theory has grown into a unifying framework that cuts across physics, biology, computer science, and the social sciences. This chapter surveys the intellectual landscape and establishes the notation and conceptual vocabulary that will be used throughout the monograph.

1.1 Historical Background

The mathematical study of networks traces its lineage to Leonhard Euler's 1736 solution of the Königsberg bridge problem, often cited as the birth of graph theory. Euler showed that it was impossible to traverse all seven bridges of Königsberg exactly once, a result that inaugurated the abstract treatment of connectivity.

The field remained largely the province of pure mathematics until the mid-twentieth century, when two developments reshaped its trajectory. First, Paul Erdős and Alfréd Rényi introduced their eponymous random graph model in 1959, providing a probabilistic framework that was immediately congenial to the statistical-mechanics tradition. Second, the rise of digital computing made it possible to collect and analyze network data at unprecedented scales.

1.2 Basic Definitions

A **network** (or **graph**) $G = (V, E)$ consists of a set of *vertices* (or *nodes*) $V = \{v_1, v_2, \dots, v_N\}$ and a set of *edges* $E \subseteq V \times V$. Throughout this monograph, $N = |V|$ denotes the number of vertices and $M = |E|$ the number of edges.

Definition 1.1 (Adjacency Matrix). The **adjacency matrix** of a simple graph on N vertices is the $N \times N$ matrix with entries

$$A_{ij} = \begin{cases} 1 & \text{if } (i, j) \in E, \\ 0 & \text{otherwise.} \end{cases}$$

For undirected graphs, A is symmetric: $A_{ij} = A_{ji}$.

For an access
overview, see
(2018), *Netw*
Oxford Univ

Euler's proof
of early topo
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networks.

The degree k_i of vertex i is the number of edges incident to it:

$$k_i = \sum_{j=1}^N A_{ij}.$$

The **degree distribution** $P(k)$ — the probability that a randomly chosen vertex has degree k — is one of the most fundamental characterizations of a network ensemble.

1.2.1 Degree Correlations and Assortativity

Networks are not fully described by their degree distribution alone. The **joint degree distribution** $P(k, k')$, which gives the probability that an edge chosen uniformly at random connects vertices of degrees k and k' , captures two-point correlations. A particularly useful scalar summary is the **assortativity coefficient** r , defined as the Pearson correlation coefficient between the degrees of adjacent vertices .

1.2.2 Clustering and Transitivity

The **clustering coefficient** C_i of vertex i measures the density of connections among its neighbors:

$$C_i = \frac{2 T_i}{k_i(k_i - 1)},$$

where T_i is the number of triangles that include vertex i . The network-wide average $\langle C \rangle$ is a standard measure of local structure.

Example 1.1 (Social vs. Biological Networks). Social networks typically exhibit high clustering ($\langle C \rangle \approx 0.3$ to 0.5), reflecting the tendency of “friends of friends” to become friends. In contrast, biological regulatory networks often show low clustering but high modularity, with dense connections within functional modules and sparse connections between them.

Chapter 2

Statistical Mechanics of Graph Ensembles

The central insight that animates this monograph is that a network ensemble can be treated as a **statistical-mechanical system**. Just as the microcanonical ensemble describes an isolated gas in terms of its accessible microstates, a **graph ensemble** describes the set of all graphs that satisfy specified macroscopic constraints.

2.1 The Microcanonical Ensemble of Graphs

Consider the set $\mathcal{G}_{N,M}$ of all simple graphs with N vertices and exactly M edges. The total number of such graphs is

$$|\mathcal{G}_{N,M}| = \binom{\binom{N}{2}}{M}.$$

In the microcanonical formulation, each graph $G \in \mathcal{G}_{N,M}$ is assigned equal probability:

$$P(G) = \frac{1}{|\mathcal{G}_{N,M}|}.$$

This is precisely the Erdős–Rényi model $G(N, M)$.

The equivalence, in the thermodynamic limit, of the microcanonical ensemble $G(N, M)$ and the canonical ensemble $G(N, p)$ with $p = M/\binom{N}{2}$ is a direct analogue of the equivalence of ensembles in classical statistical mechanics.

2.2 The Canonical Ensemble: Exponential Random Graphs

The canonical formulation is considerably more flexible. Given a set of graph observables $\{X_\mu(G)\}_{\mu=1}^K$ and their prescribed expected values $\{\langle X_\mu \rangle\}$, the maximum-entropy distribution

The analogy
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mechanics.

takes the exponential form

$$P(G) = \frac{1}{Z} \exp\left(-\sum_{\mu=1}^K \theta_{\mu} X_{\mu}(G)\right),$$

where $Z = \sum_G \exp(-\sum_{\mu} \theta_{\mu} X_{\mu}(G))$ is the partition function and the Lagrange multipliers $\{\theta_{\mu}\}$ are determined by the constraint equations $\langle X_{\mu} \rangle = -\partial \log Z / \partial \theta_{\mu}$.

Definition 2.1 (Hamiltonian of a Graph Ensemble). The quantity

$$\mathcal{H}(G) = \sum_{\mu=1}^K \theta_{\mu} X_{\mu}(G)$$

is called the **graph Hamiltonian**. It assigns an “energy” to each graph, with lower-energy graphs being more probable under the ensemble.

2.2.1 The Configuration Model

The most important special case is the configuration model, where the prescribed observables are the degree sequence $\{k_i\}_{i=1}^N$. The Hamiltonian becomes

$$\mathcal{H}(G) = \sum_{i=1}^N \theta_i k_i(G) = \sum_{i<j} (\theta_i + \theta_j) A_{ij},$$

and the probability factorizes over edge variables:

$$P(G) = \prod_{i<j} p_{ij}^{A_{ij}} (1 - p_{ij})^{1-A_{ij}}, \quad p_{ij} = \frac{e^{-(\theta_i + \theta_j)}}{1 + e^{-(\theta_i + \theta_j)}}.$$

Theorem 2.1 (Existence of the MaxEnt Ensemble). *For any sequence of real numbers $\{\langle k_i \rangle\}_{i=1}^N$ satisfying $0 < \langle k_i \rangle < N - 1$ for all i , there exists a unique set of Lagrange multipliers $\{\theta_i\}$ such that the canonical ensemble reproduces the prescribed expected degrees.*

Proof. The log-likelihood $\mathcal{L}(\theta) = -\log Z + \sum_i \theta_i \langle k_i \rangle$ is strictly convex in θ , since its Hessian $\partial^2 \mathcal{L} / \partial \theta_i \partial \theta_j = \text{Cov}(A_{ij})$ is positive-definite on the interior of the feasible region. Its gradient $\nabla_{\theta_i} \mathcal{L} = \langle k_i \rangle - \sum_{j \neq i} p_{ij}$ therefore vanishes at a unique point, and the map $\theta \mapsto \{\langle k_i \rangle\}$ is a diffeomorphism onto the space of feasible degree sequences. $\square$

2.3 Phase Transitions in Network Ensembles

Theorem 2.2 (Percolation Transition). *In the Erdős–Rényi ensemble $G(N, p)$ with $p = c/N$, the expected size of the largest connected component undergoes a second-order phase transition at $c = 1$. For $c < 1$, the largest component has size $O(\log N)$; for $c > 1$, a **giant component** emerges that contains a finite fraction of all vertices.*

The analogy with equilibrium critical phenomena runs deep. The order parameter $m(c) = S/N$,

where S is the size of the giant component, scales as $m \sim (c - 1)^\beta$ with $\beta = 1$ near the critical point — precisely the mean-field exponent .

2.3.1 Beyond the Configuration Model

While the configuration model captures arbitrary degree distributions, it ignores higher-order correlations. The **stochastic block model** (SBM) introduces community structure by partitioning vertices into Q blocks and assigning connection probabilities p_{qr} between blocks q and r . The SBM Hamiltonian is

$$\mathcal{H}(G) = \sum_{q,r=1}^Q \theta_{qr} e_{qr}(G),$$

where e_{qr} counts edges between blocks q and r .

Remark 2.1 (Identifiability). A recurring theme in the statistical mechanics of networks is the tension between model richness and parameter identifiability. The SBM with $O(Q^2)$ free parameters is unidentifiable when the block structure is unknown, motivating the extensive literature on community detection and model selection in networks.

Corollary 2.3. *Under mild regularity conditions, the maximum-likelihood estimate of the SBM parameters is consistent as $N \rightarrow \infty$ provided that the average degree diverges no faster than $\log N$.*

This is not a
Erdős–Rényi
a fully conn
geometry, an
theory is exa

Part II

Applications and Case Studies

All models are wrong, but some are useful.

George E. P. Box

Color Plates

Selected illustrations referenced throughout the text.

Color versions are available in the electronic edition.

Plate 1

Power-law degree distributions. Empirical degree distributions for three real-world networks, plotted on double-logarithmic axes. Left: the Internet at the autonomous-system level ($N \approx 25,000$, estimated exponent $\gamma = 2.1$). Center: a co-authorship network in condensed-matter physics ($N \approx 40,000$, $\gamma = 2.4$). Right: word co-occurrence in *Moby Dick* ($N \approx 17,000$, $\gamma = 2.8$). The dashed lines show maximum-likelihood power-law fits; shaded regions indicate 95% confidence intervals obtained via parametric bootstrap with $B = 10,000$ resamples.

Plate 2

Phase diagram of the correlated Erdős–Rényi ensemble. The axes represent the average degree c and the assortativity strength α . Region I (blue): subcritical regime with no giant component and exponentially decaying component-size distribution. Region II (green): a single giant component spanning approximately $m(c) \cdot N$ vertices. Region III (orange): a *bimodal* phase in which two giant components coexist, stabilized by strong degree–degree correlations. The critical line separating Regions I and II is second-order; the transition between II and III is first-order. The tricritical point at $(c, \alpha) \approx (2.4, 0.62)$ marks the meeting of the two phase boundaries.

Plate 3

Community detection via belief propagation. Application of the degree-corrected stochastic block model to a friendship network of $N = 1,034$ students at a U.S. high school. Left panel: the adjacency matrix reordered according to the inferred block assignment, revealing five distinct communities that correspond closely to self-reported social groups. Right panel: the same matrix randomized while preserving the degree sequence, demonstrating that the block structure is statistically significant ($p < 10^{-4}$ under a parametric bootstrap test). Colors indicate edge density within and between blocks, normalized by the expected density under the configuration-model null.

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