

Basin Percolation in the LSR Associative Memory

1 Setup: LSR energy and the hard wall

The Log-Sum-ReLU (LSR) associative memory with M patterns ξ^1, \dots, ξ^M on the sphere $S^{N-1}(\sqrt{N})$ has energy

$$E(\mathbf{S}) = -\frac{N}{b} \ln \sum_{\mu=1}^M [1 - b + b \varphi_\mu(\mathbf{S})]_+, \quad b = 2 + \sqrt{2}, \quad (1)$$

where $\varphi_\mu(\mathbf{S}) = (\xi^\mu \cdot \mathbf{S})/N$ is the overlap and $[\cdot]_+ = \max(0, \cdot)$.

The hard wall activates at $\varphi_c = (b-1)/b = 1/\sqrt{2} \approx 0.707$: pattern μ contributes to the energy only when $\varphi_\mu(\mathbf{S}) > \varphi_c$. When no pattern is active, the argument of the logarithm vanishes and $E = +\infty$.

Basin of a single pattern. For an isolated target ξ^1 , the accessible region is the spherical cap

$$X_0 = \{ \mathbf{S} \in S^{N-1}(\sqrt{N}) : \varphi_1(\mathbf{S}) > \varphi_c \}, \quad (2)$$

with an infinite energy barrier at the boundary $\varphi_1 = \varphi_c$. The state is permanently confined to X_0 .

2 Joint basins and continuous connectivity

With M patterns, the accessible region is the set where $\sum_\mu [\cdots]_+ > 0$, i.e. where at least one pattern is active. Two individual basins X_μ and X_ν form a *joint basin* if their intersection $X_\mu \cap X_\nu \neq \emptyset$.

Along any continuous path through $X_\mu \cap X_\nu$, both terms are simultaneously positive, so the energy is always finite—the state never encounters $\ln 0$. The energy does increase at the “bottleneck” (narrowest part of the overlap corridor), creating a finite barrier

$$\Delta E \sim \frac{N}{2b} \ln \frac{1}{q}, \quad (3)$$

where $q = (\xi^\mu \cdot \xi^\nu)/N$ is the mutual overlap, but this barrier is finite for any $q > 0$.

Geometric overlap criterion. Two spherical caps of angular radius $\theta_c = \arccos \varphi_c = \pi/4$ overlap when the angle between their centers satisfies $\theta < 2\theta_c = \pi/2$, i.e. when

$$q = \frac{\xi^\mu \cdot \xi^\nu}{N} > \cos\left(\frac{\pi}{2}\right) = 0. \quad (4)$$

Since random patterns have $q \sim \mathcal{N}(0, 1/N)$, approximately $M/2$ patterns satisfy $q > 0$ geometrically. However, the energy barrier (??) scales as $\sim (N/4b) \ln N$ for $q \sim 1/\sqrt{N}$, making such corridors thermally inaccessible.

3 Percolation framework

3.1 Thermal connectivity threshold

We define two basins as *thermally connected* if the energy barrier along the optimal path through their overlap is comparable to the thermal energy scale T . This requires a minimum mutual overlap $q > q_{\text{eff}}(T)$.

At zero temperature, only basins with a wide, low-barrier corridor are connected. The natural hard-wall threshold is $q_{\text{eff}} = \varphi_c$, since patterns with $q > \varphi_c$ are already active at the target center (the term $[1 - b + b q]_+ > 0$).

3.2 Poisson model for the number of neighbors

For random patterns on $S^{N-1}(\sqrt{N})$, the overlap $q = (\boldsymbol{\xi}^\mu \cdot \boldsymbol{\xi}^1)/N$ with a fixed target is approximately $\mathcal{N}(0, 1/N)$ for large N . The number of patterns with $q > \varphi_c$ follows a Poisson distribution with rate

$$\lambda = M \cdot \Pr(q > \varphi_c) \approx \frac{\exp(N(\alpha - \varphi_c^2/2))}{\varphi_c \sqrt{2\pi N}}, \quad (5)$$

where $M = e^{N\alpha}$ and we used the Mill's ratio approximation for the Gaussian tail.

Finite- N correction. The exact overlap distribution on S^{N-1} has density $f(q) \propto (1-q^2)^{(N-3)/2}$, which has a lighter tail than the Gaussian for moderate N . The Mill's ratio therefore *overestimates* λ ; the true percolation threshold is shifted to higher α for finite N . This correction vanishes as $N \rightarrow \infty$.

3.3 Percolation transition

The basins of M random patterns form a random geometric graph on S^{N-1} : each basin is a node, and two nodes are connected if their mutual overlap exceeds q_{eff} . The connected component containing the target undergoes a **percolation transition** at $\lambda = 1$.

Setting $\lambda = 1$ in (??) with $q_{\text{eff}} = \varphi_c$:

$$\boxed{\alpha_c = \frac{\varphi_c^2}{2} + \frac{\ln(\varphi_c \sqrt{2\pi N})}{N} \xrightarrow{N \rightarrow \infty} \frac{\varphi_c^2}{2} = \frac{1}{4} = 0.25.} \quad (6)$$

This coincides exactly with the zero-temperature critical capacity $\alpha_{\text{th}} = \frac{1}{2}(1 - 1/b)^2 = \frac{1}{2}\varphi_c^2$ from the mean-field theory.

Physical picture.

- **Subcritical** ($\alpha < \alpha_c$): $\lambda < 1$; the target's connected cluster has $O(1)$ patterns. The state is confined near the target \rightarrow retrieval.
- **Critical** ($\alpha \approx \alpha_c$): $\lambda \approx 1$; cluster size diverges; critical fluctuations.
- **Supercritical** ($\alpha > \alpha_c$): $\lambda > 1$; a giant connected component forms. The state delocalizes among exponentially many basins \rightarrow paramagnetic phase.

3.4 Temperature dependence

At temperature $T > 0$, thermal fluctuations lower the effective barrier, so narrower corridors become traversable: $q_{\text{eff}}(T) < \varphi_c$. The percolation threshold shifts to

$$\alpha_c(T) \approx \frac{q_{\text{eff}}(T)^2}{2}, \quad (7)$$

which decreases with T —the retrieval region shrinks, consistent with the phase diagram.

From the mean-field free energy calculation: $\alpha_c(T) = \frac{1}{2}[1 - f_{\text{ret}}(T)]^2$, where $f_{\text{ret}}(T) = u(\varphi(T)) - T s(\varphi(T))$ is the retrieval free energy. Identifying $q_{\text{eff}}(T) = 1 - f_{\text{ret}}(T)$ connects the percolation threshold to the thermodynamic phase boundary.

4 BFS depth: chain connections

Starting from the target, we explore the basin connectivity graph via breadth-first search:

1. **Depth 0:** target pattern ξ^1 with basin X_0 .
2. **Depth 1:** all patterns μ with $(\xi^\mu \cdot \xi^1)/N > \varphi_c$ ($K_1 \sim \text{Poisson}(\lambda)$ patterns).
3. **Depth 2:** for each depth-1 pattern μ , find patterns ν with $(\xi^\mu \cdot \xi^\nu)/N > \varphi_c$ that are *not* already at depth 1.

Depth-2 patterns are absent for large N . For a depth-1 pattern μ (overlap $\varphi_\mu > \varphi_c$ with target) and a candidate depth-2 pattern ν (overlap $\varphi_\nu < \varphi_c$ with target), their mutual overlap is

$$\frac{\xi^\mu \cdot \xi^\nu}{N} \approx \varphi_\mu \cdot \varphi_\nu + \sqrt{(1 - \varphi_\mu^2)(1 - \varphi_\nu^2)} \frac{u_\mu \cdot u_\nu}{\sqrt{N-1}}, \quad (8)$$

where u_μ, u_ν are random unit vectors in the $(N-1)$ -dimensional subspace perpendicular to the target.

The deterministic part satisfies $\varphi_\mu \cdot \varphi_\nu < \varphi_c^2 = 1/2 < \varphi_c = 1/\sqrt{2}$, and the random fluctuation $\sim \mathcal{O}(1/\sqrt{N})$ cannot bridge the gap $\varphi_c - \varphi_c^2 \approx 0.207$ for large N .

At finite N , the fluctuation has standard deviation $\sim 0.5/\sqrt{N-1}$, requiring a $\sim 0.207\sqrt{N}$ sigma event for a depth-2 connection. For $N = 25$: $\sim 1\sigma$ (rare but possible); for $N = 50$: $\sim 1.5\sigma$ (very rare); for $N \geq 100$: negligible.

5 Simulation design

The accompanying script `percolation_LSR.jl` tests four aspects of the percolation hypothesis:

1. **Panel 1: Analytical** $\lambda(\alpha)$ for $N = 30, 50, 75, 150$. All curves cross $\lambda = 1$ near $\alpha \approx 0.25$, with finite- N corrections $\sim \ln N/N$ shifting the threshold to higher α .
2. **Panel 2: Direct neighbor counting** ($N = 25, 3000$ realizations). Generate $M = e^{N\alpha}$ random patterns on $S^{N-1}(\sqrt{N})$, count neighbors K with overlap $> \varphi_c$. Compare histogram with Poisson(λ). The Mill's ratio approximation overestimates λ at moderate N because the exact spherical tail is lighter than the Gaussian.

3. **Panel 3: Branching process survival** ($N = 50$). Simulate a Galton–Watson process with Poisson(λ) offspring. The survival probability transitions from 0 to 1 at α_c , sharpening with BFS depth (analogous to increasing system size in percolation).
4. **Panel 4: BFS depth analysis** ($N = 25, 200$ realizations). Count depth-1 neighbors K_1 and new depth-2 neighbors K_2 . Confirms that $\langle K_2 \rangle \ll \langle K_1 \rangle$: the BFS effectively terminates at depth 1.

6 Implications for the simulation protocol

The percolation picture justifies the Poisson pattern reduction scheme (v5 approach) with an important correction: **the background energy S_{bg} should be removed**.

1. The energy computed from just the $K + 1$ retained patterns (target + K active neighbors) is *exact* for all states reachable from the target basin. Patterns outside the connected component are separated by infinite barriers and cannot be reached.
2. The S_{bg} term in v5 artificially smooths the energy landscape, removing the infinite barriers that confine the state. This distorts the dynamics near the phase boundary.
3. With the Poisson scheme (no S_{bg}), memory scales as $\mathcal{O}(N \cdot K_{\max})$ instead of $\mathcal{O}(N \cdot M)$, enabling much larger N for sharper phase boundaries.
4. For $\alpha > \alpha_c$ (supercritical), K grows exponentially but can be capped at some K_{cap} without affecting the measured overlap φ at the target—the state delocalizes among the K_{cap} nearest basins, which is sufficient to show non-retrieval.

Connection to the Hopfield model. In the classical Hopfield model, the coupling structure creates an exponentially large number of spin-glass metastable states via random-matrix frustration. In the LSR model, the log- \sum -exp energy function confines the state to discrete basins separated by infinite barriers. The phase transition is purely *geometric*—basin percolation—rather than arising from frustrated interactions. This explains the absence of the spin-glass phase in dense associative memories.

References

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