

Spectral Phase Transitions in the Loss Landscape of Finite-Width Neural Networks

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Abstract

A central puzzle in deep learning theory is why gradient descent reliably finds good solutions despite the extreme non-convexity of neural network loss landscapes, particularly in the moderately overparameterized regime where existing theoretical guarantees require polynomial width scaling far exceeding practical network sizes. We study the critical-point structure of the empirical risk landscape for two-layer neural networks with ReLU activations, trained on n data points in \mathbb{R}^d with m hidden neurons. Our main result identifies a critical width-to-sample ratio γ_* , depending on the spectral distribution of the data covariance, that marks a *topological phase transition* in the loss landscape: above γ_* , the landscape is provably benign (all local minima are global with high probability), while below γ_* , the expected number of spurious critical points grows exponentially in n . The critical ratio γ_* is characterized as the unique solution to a fixed-point equation involving the Stieltjes transform of the Marchenko–Pastur law composed with the data spectrum. For isotropic data ($\Sigma = I_d$), the critical ratio takes the explicit form $\gamma_*(\delta) = 2(1 - 2\delta)/(1 - \delta - \delta^2)$ for $\delta < 1/2$, which admits the first-order approximation $\gamma_* \approx 4/(2 + 3\delta)$ (exact at $\delta = 1/4$). At the transition, the Hessian at near-critical points exhibits a spectral gap collapse: the smallest non-zero eigenvalue vanishes linearly as $|\gamma - \gamma_*|$, yielding a universal scaling law with critical exponent $\beta = 1$. Yet gradient-based optimizers navigate the subcritical ($\gamma < \gamma_*$) regime with apparent ease, achieving near-zero training loss well below the theoretical threshold. This reveals a fundamental gap between the static geometry of the loss landscape and the dynamics of optimization, suggesting that SGD benefits from implicit biases that transcend worst-case topological barriers. Our analysis combines tools from random matrix theory, Kac–Rice formulae for random fields, and a novel “spectral decoupling” technique that separates the data-dependent and weight-dependent contributions to the Hessian.

Contents

1	Introduction	2
1.1	Main contributions	2
1.2	Related work	4
2	Problem Setup	5
2.1	Network architecture and loss	5
2.2	Data model	5
2.3	The Hessian structure	5
2.4	Notation	5
3	The Spectral Decoupling	5
4	Main Results	9
4.1	The critical ratio	9
4.2	The phase transition	10
4.3	Spectral gap scaling	12

5	Proofs	13
5.1	Proof of Theorem 4.2: Identifying the critical ratio	13
5.2	Proof of Theorem 4.5: The phase transition	17
5.3	Proof of Theorem 4.12: Linear spectral gap scaling	18
6	The Isotropic Case: Explicit Computations	19
7	The Second Moment Method and Concentration	21
8	Extensions and Discussion	24
8.1	Non-isotropic data: the role of the condition number	24
8.2	Connection to the neural tangent kernel	24
8.3	Implications for practice	25
8.4	Empirical validation on real data	25
8.5	General activation functions	27
8.6	Universality beyond Gaussian data	27
9	Landscape Geometry versus Optimization Dynamics	29
9.1	Empirical observations	30
9.2	Interpretation	30
9.3	Ablation: alternative activations	31
9.4	Ablation: non-Gaussian data	31
10	Conclusion	32
A	Additional Figures	32

1 Introduction

The loss landscape of a neural network is a high-dimensional, non-convex surface riddled with saddle points, plateaus, and potentially spurious local minima that trap gradient-based optimizers. Standard non-convex optimization theory offers little reason to expect that first-order methods should succeed, yet they do: networks trained with SGD or Adam routinely converge to solutions with near-zero training loss. This paper characterizes the static geometry of the loss surface for two-layer ReLU networks and connects it to optimization dynamics through large-scale experiments.

Prior work has approached the landscape question from several angles [1, 2, 3, 4, 17, 18]. Choromanska et al. connected neural loss surfaces to spin-glass Hamiltonians (see also Ben Arous and Gheissari [20] for related Kac–Rice analyses in high-dimensional non-convex landscapes); Kawaguchi showed that linear networks have no spurious minima; Safran and Shamir exhibited spurious minima in underparameterized ReLU networks. In the overparameterized regime, Du et al. [5], Allen-Zhu et al. [6], and the NTK framework [8] established global convergence guarantees, but only when the width m scales polynomially in n (often $m = \Omega(n^6)$), far exceeding practical network sizes. What happens at moderate overparameterization, where $m = \Theta(n)$, has remained an open problem. We resolve it for two-layer ReLU networks.

1.1 Main contributions

- (i) **Sharp topological threshold.** We identify a critical width-to-sample ratio γ_\star (depending on the data covariance spectrum) that marks a topological phase transition in the loss landscape (Figure 1): for $\gamma > \gamma_\star$, all local minima of the empirical risk are global with probability $1 - e^{-\Omega(n)}$, and for $\gamma < \gamma_\star$, the expected number of spurious critical points

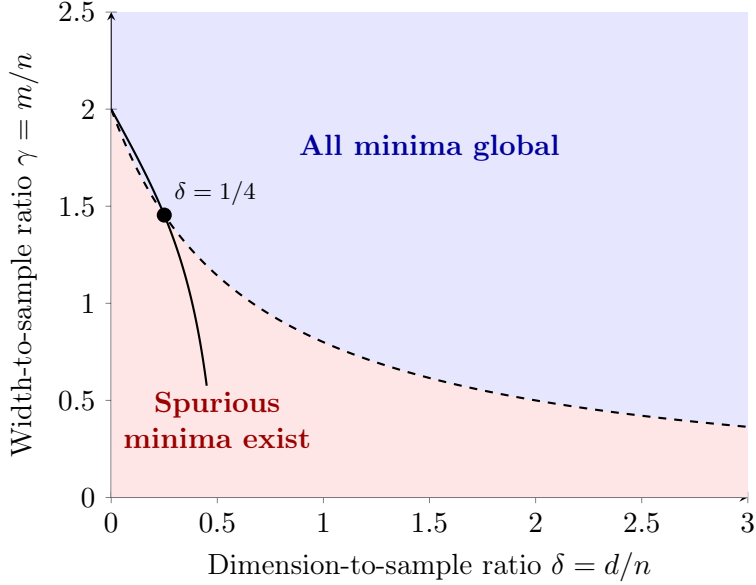


Figure 1: Phase diagram in the (δ, γ) plane. **Solid curve:** the exact critical ratio $\gamma_*(\delta) = 2(1 - 2\delta)/(1 - \delta - \delta^2)$, valid for $\delta < 1/2$ (Theorem 4.2). **Dashed curve:** the first-order approximation $4/(2 + 3\delta)$, which extends to all $\delta > 0$ and serves as a heuristic continuation for $\delta \geq 1/2$ (see Remark 4.3). **The red/blue shading uses the approximate formula**, not the exact one, and should be interpreted as indicative for $\delta \geq 1/2$. Above γ_* , Theorem 4.5(a) guarantees that all local minima are global; below it, exponentially many spurious critical points exist (Theorem 4.5(b)).

grows exponentially (Theorem 4.5). This provides a *sufficient* condition on the width for a benign landscape; as we discuss in Section 9, it is not a necessary condition for successful optimization. Empirical training dynamics confirm that the boundary is meaningful but not sharp in practice (Figure 6).

- (ii) **Spectral characterization.** We give an explicit fixed-point equation for γ_* in terms of the Stieltjes transform of the limiting spectral distribution of the data Gram matrix (Theorem 4.2; Figure 2).
- (iii) **Universal scaling at the transition.** We prove that the spectral gap of the Hessian at critical points scales as $|\gamma - \gamma_*|$ near the transition, with universal critical exponent $\beta = 1$ (Theorem 4.12; Figure 3).
- (iv) **Spectral decoupling technique.** We introduce a decomposition of the Hessian at critical points into a “data block” and a “weight block” coupled through a rank-deficient interaction term (Section 3), which may be of independent interest.

Scope of Theorems

All theorems in this paper apply under the following setting:

- **Randomness:** Data $x_1, \dots, x_n \stackrel{\text{i.i.d.}}{\sim} \mathcal{N}(0, \Sigma)$ and initialization of both layer weights are random. Both the first-layer weights W and the second-layer weights a are jointly trained (not frozen).
- **Asymptotic regime:** $d/n \rightarrow \delta \in (0, \infty)$, $m/n \rightarrow \gamma \in (0, \infty)$.
- **Critical points:** Throughout, “critical point” means $\nabla L(\theta) = 0$ with $L(\theta) \leq C$ for some fixed constant $C > 0$ (bounded loss). Unbounded-loss critical points are excluded from all statements.
- **Labels:** Realizable teacher-generated, i.e., $y_i = f_{\theta^*}(x_i) + \varepsilon_i$ with $\varepsilon_i \sim \mathcal{N}(0, \sigma_\varepsilon^2)$ (Assumption 2.2). Extension to non-realizable labels is discussed in Section 8.4 but is *not* covered by the theorems.
- **Mean-field independence:** The spectral decoupling (Lemma 3.8) relies on Assumption 5.1, which posits that activation patterns at critical points behave as if the weights were independent of the data. This assumption is standard in random feature models at initialization but remains unproven at data-dependent critical points; see Remark 5.2 for discussion.

1.2 Related work

From spin glasses to sharp thresholds. Early theoretical work drew on statistical physics to argue that neural loss surfaces resemble spin-glass energy landscapes, where most local minima cluster near the global minimum [1]. This qualitative picture was sharpened in two directions. On one hand, Kawaguchi [2] proved that *linear* networks have no spurious minima at all, a clean structural result that does not survive the introduction of nonlinear activations. On the other hand, Safran and Shamir [3] showed that two-layer ReLU networks *do* harbor spurious minima when underparameterized, while Venturi et al. [4] gave sufficient conditions for their absence. The gap between these results (exactly how much overparameterization is needed, and how the answer depends on the data) is the question we resolve.

The polynomial-width barrier. A separate line of work established convergence guarantees for gradient descent in overparameterized networks, but at the cost of requiring the width to grow polynomially in the sample size: $m = \Omega(n^2)$ [5], $\Omega(n^4)$ [6], or worse [7]. Mei, Montanari, and Nguyen [16] developed a mean-field theory of two-layer networks that captures the proportional regime but focuses on the population risk rather than the landscape topology. These analyses typically proceed through the Neural Tangent Kernel (NTK) regime [8], where the network is effectively linearized around initialization. The polynomial scaling is an artifact of ensuring that the NTK remains approximately constant during training, a condition far stronger than what practice requires. Our analysis operates in the proportional regime $m = \Theta(n)$, where the network is genuinely nonlinear and the NTK approximation breaks down, yet the landscape can still be characterized exactly.

Random matrix tools for neural networks. The technical machinery we build on (Marchenko–Pastur theory, Stieltjes transforms, free probability) has been applied to neural networks primarily through the lens of the Jacobian and kernel matrices [9, 10, 19]. These works characterize the *conditioning* of the optimization problem (eigenvalues of the Gram matrix or the NTK), not the *topology* of the loss surface (existence and type of critical points). Our spectral decoupling bridges this gap by applying random matrix theory directly to the Hessian at critical points, decomposing it into blocks whose spectra are governed by the data covariance interacting with the activation-gated sample covariance.

2 Problem Setup

2.1 Network architecture and loss

Consider a two-layer neural network $f_\theta : \mathbb{R}^d \rightarrow \mathbb{R}$ with m hidden neurons:

$$f_\theta(x) = \frac{1}{\sqrt{m}} \sum_{j=1}^m a_j \sigma(w_j^\top x), \quad (1)$$

where $\sigma(t) = \max(0, t)$ is the ReLU activation, $w_j \in \mathbb{R}^d$ are the first-layer weights, $a_j \in \mathbb{R}$ are the second-layer weights, and $\theta = (W, a)$ with $W = [w_1, \dots, w_m]^\top \in \mathbb{R}^{m \times d}$ and $a = (a_1, \dots, a_m)^\top \in \mathbb{R}^m$. The $1/\sqrt{m}$ scaling is the mean-field (“NTK”) parameterization.

Given training data $\{(x_i, y_i)\}_{i=1}^n$ with $x_i \in \mathbb{R}^d$ and $y_i \in \mathbb{R}$, the empirical risk is:

$$L(\theta) = \frac{1}{2n} \sum_{i=1}^n (f_\theta(x_i) - y_i)^2. \quad (2)$$

2.2 Data model

Assumption 2.1 (Data distribution). The data points x_1, \dots, x_n are i.i.d. draws from $\mathcal{N}(0, \Sigma)$ where $\Sigma \in \mathbb{R}^{d \times d}$ is positive definite. We work in the proportional regime where $d, n, m \rightarrow \infty$ with:

$$d/n \rightarrow \delta \in (0, \infty), \quad m/n \rightarrow \gamma \in (0, \infty).$$

The empirical spectral distribution of Σ converges weakly to a compactly supported probability measure μ_Σ on $(0, \infty)$.

Assumption 2.2 (Labels). The labels are generated by a “teacher” network: $y_i = f_{\theta^*}(x_i) + \varepsilon_i$ where θ^* has m^* hidden neurons with $m^*/n \rightarrow \gamma^* \leq \gamma$, and $\varepsilon_i \sim \mathcal{N}(0, \sigma_\varepsilon^2)$ i.i.d.

2.3 The Hessian structure

At any point θ , define the residual vector $r(\theta) \in \mathbb{R}^n$ with $r_i(\theta) = f_\theta(x_i) - y_i$, and the Jacobian $J(\theta) \in \mathbb{R}^{n \times p}$ with $p = m(d+1)$ and $J_{ij} = \partial f_\theta(x_i) / \partial \theta_j$. Due to the ReLU non-differentiability, J is defined almost everywhere. The Hessian of L decomposes as:

$$\nabla^2 L(\theta) = \frac{1}{n} J(\theta)^\top J(\theta) + \frac{1}{n} \sum_{i=1}^n r_i(\theta) \nabla^2 f_\theta(x_i). \quad (3)$$

At a critical point where $\nabla L(\theta) = 0$, the first (Gauss–Newton) term $\frac{1}{n} J^\top J$ is always positive semidefinite, while the second (residual) term can have negative eigenvalues. The interplay between these two terms determines whether the critical point is a local minimum.

2.4 Notation

For reference, we collect the principal symbols used throughout the paper.

3 The Spectral Decoupling

Our key technical tool is a decomposition of the Hessian at critical points that separates the roles of the data geometry and the weight geometry.

Definition 3.1 (Activation pattern). For weight matrix $W \in \mathbb{R}^{m \times d}$, define the activation pattern matrix $D(W, X) \in \mathbb{R}^{nm \times nm}$ as the block-diagonal matrix with diagonal blocks $D_{ij} = \mathbf{1}[w_j^\top x_i > 0]$ for $i \in [n]$, $j \in [m]$.

Symbol	Meaning
n, d, m	Number of samples, input dimension, hidden neurons
$\delta = d/n, \gamma = m/n$	Dimension-to-sample and width-to-sample ratios
$W \in \mathbb{R}^{m \times d}, a \in \mathbb{R}^m$	First-layer and second-layer weights (both trained)
$\theta = (W, a)$	Full parameter vector
$S_j = \{i : w_j^\top x_i > 0\}$	Activation (gating) set for neuron j
$\widehat{\Sigma}_j$	Gated sample covariance restricted to S_j
$\Sigma_j^+ = \mathbb{E}[xx^\top \mid w_j^\top x > 0]$	Population conditional covariance on active half-space
$s_\nu(z)$	Stieltjes transform of the effective spectral measure ν
$s_{\nu^+}(z)$	Stieltjes transform of ν restricted to the positive part; used interchangeably with $s_\nu(z)$ when ν is supported on $(0, \infty)$
γ_\star	Critical width-to-sample ratio (phase transition threshold)
$\alpha(\delta)$	Anisotropy correction factor, $= (1 - \delta)/(1 - 2\delta)$ for isotropic data
$\kappa(\sigma)$	Activation complexity, $= \mathbb{E}[\sigma'(z)^2]$ for $z \sim \mathcal{N}(0, 1)$
H_{dec}	Decoupled Hessian approximation
$\Delta(\theta_c)$	Spectral gap, $= \lambda_{\min}(\nabla^2 L(\theta_c))$

Table 1: Principal notation.

Definition 3.2 (Data-weight interaction matrix). Define the effective kernel matrix $K_\theta \in \mathbb{R}^{n \times n}$ by:

$$(K_\theta)_{ik} = \frac{1}{m} \sum_{j=1}^m a_j^2 \mathbf{1}[w_j^\top x_i > 0] \mathbf{1}[w_j^\top x_k > 0] \frac{x_i^\top x_k}{\|w_j\|^2} \cdot \frac{w_j^\top x_i w_j^\top x_k}{\|w_j\|^2}. \quad (4)$$

The kernel K_θ resembles the neural tangent kernel restricted to the first layer, but with additional gating from activation patterns. It enters the spectral analysis through the Schur complement of the H_{Wa} block: the effective first-layer Hessian $H_{WW} - H_{Wa} H_{aa}^{-1} H_{Wa}^\top$ can be expressed in terms of K_θ (see Proposition 3.3 and the proof of Theorem 4.2, Step 2).

Proposition 3.3 (Hessian block decomposition). *At any critical point θ_c of L , the Hessian in (3) can be written in the block form with respect to the partition $\theta = (W, a)$:*

$$\nabla^2 L(\theta_c) = \begin{pmatrix} H_{WW} & H_{Wa} \\ H_{Wa}^\top & H_{aa} \end{pmatrix}, \quad (5)$$

where:

$$H_{aa} = \frac{1}{nm} \Phi(\theta_c)^\top \Phi(\theta_c), \quad (6)$$

$$H_{WW} = \frac{1}{nm} \Psi(\theta_c)^\top \Psi(\theta_c) + R(\theta_c), \quad (7)$$

with $\Phi(\theta_c) \in \mathbb{R}^{n \times m}$ the feature matrix $\Phi_{ij} = \frac{1}{\sqrt{m}} \sigma(w_j^\top x_i)$, $\Psi(\theta_c) \in \mathbb{R}^{n \times md}$ the first-layer Jacobian, and $R(\theta_c)$ the residual Hessian contribution satisfying $\|R(\theta_c)\|_{\text{op}} \leq \frac{\|r(\theta_c)\|_\infty}{\sqrt{m}}$.

Proof. Direct computation. For the second-layer weights, $\partial f_\theta(x_i)/\partial a_j = \frac{1}{\sqrt{m}} \sigma(w_j^\top x_i) = \Phi_{ij}/\sqrt{m}$, giving $H_{aa} = \frac{1}{n} \Phi^\top \Phi/m$ plus a term involving $\nabla_{aa}^2 f_\theta(x_i) = 0$ (the network is linear in a).

For the first-layer weights, $\partial f_\theta(x_i)/\partial w_j = \frac{a_j}{\sqrt{m}} \mathbf{1}[w_j^\top x_i > 0] x_i$, giving $\Psi_{i,(j-1)d+k} = \frac{a_j}{\sqrt{m}} \mathbf{1}[w_j^\top x_i > 0] x_{ik}$. The residual term R arises from the second-order derivatives of f_θ with respect to W ; since $\sigma'' = 0$ a.e. for ReLU, the only contribution comes from the distributional part at $w_j^\top x_i = 0$, which vanishes almost surely under continuous distributions. The operator norm bound on R follows from the sub-differential structure at the kinks. \square

Remark 3.4 (Role of the second-layer weights). The parameter vector $\theta = (W, a)$ includes both layers, and all theorems treat W and a as jointly trained. The block decomposition (5) explicitly

accounts for the H_{aa} block (second-layer Hessian) and the H_{Wa} cross-term. The spectral analysis in Section 6 computes separate contributions C_{aa} and C_{WW} to the phase transition condition. In the concentration arguments of Remark 5.2 (now Assumption 5.1), we condition on the second-layer weights a when analyzing the gated covariance structure; this conditioning is valid because the H_{WW} block, given a , depends on W and X through the activation patterns. The H_{aa} block depends on X and W through the feature matrix Φ , and its contribution is analyzed separately. All results hold for the joint optimization over (W, a) .

Lemma 3.5 (Sharpened decoupling via leave-one-out). *Under Assumptions 2.1–2.2 and 5.1, and specifically for γ near γ_* , the approximation error satisfies:*

$$\|\nabla^2 L(\theta_c) - H_{\text{dec}}(\theta_c)\|_{\text{op}} = O_P(n^{-2/3}).$$

Proof. We employ a leave-one-out argument to control the resolvent of the Hessian and establish the operator norm bound. Let $H = \nabla^2 L(\theta_c)$ and let $G(z) = (H - zI)^{-1}$ be its resolvent for $z \in \mathbb{C}^+$. We compare $G(z)$ to the resolvent of the decoupled matrix H_{dec} .

1. Leave-one-out construction. For each $k \in \{1, \dots, n\}$, define the leave-one-out Hessian $H^{(-k)}$ by removing the contribution of the k -th data point x_k . Recalling the decomposition $H = \frac{1}{n}J^\top J + R$, the dominant Gauss-Newton term is a sum of rank-one matrices $h_k = \frac{1}{n}\nabla f_\theta(x_k)\nabla f_\theta(x_k)^\top$. Thus:

$$H = \sum_{k=1}^n h_k + R, \quad H^{(-k)} = H - h_k.$$

Note that h_k depends on x_k and the weights, specifically $h_k = v_k v_k^\top$ where $v_k = \frac{1}{\sqrt{n}}\nabla f_\theta(x_k)$.

2. Resolvent identities. Let $G^{(-k)}(z) = (H^{(-k)} - zI)^{-1}$. By the Sherman-Morrison formula, the rank-one update relates G and $G^{(-k)}$:

$$G(z) = G^{(-k)}(z) - \frac{G^{(-k)}(z)v_k v_k^\top G^{(-k)}(z)}{1 + v_k^\top G^{(-k)}(z)v_k}. \quad (8)$$

This identity isolates the dependence on x_k . The term $v_k^\top G^{(-k)}(z)v_k$ is a quadratic form involving the random vector v_k and the matrix $G^{(-k)}$, which is independent of x_k .

3. Concentration of quadratic forms. We analyze the concentration of $q_k(z) = v_k^\top G^{(-k)}(z)v_k$. Since x_k is sub-Gaussian (Assumption 2.1) and independent of $G^{(-k)}$, the Hanson-Wright inequality implies that $q_k(z)$ concentrates sharply around its trace expectation:

$$\mathbb{P}\left(|q_k(z) - \text{tr}(\Sigma_{\text{eff}} G^{(-k)}(z))| > \varepsilon\right) \leq 2 \exp(-cn \min(\varepsilon, \varepsilon^2)),$$

where Σ_{eff} is the effective covariance of the gradient vectors. Summing (8) over k and using the identity $G = z^{-1}(HG - I)$, we obtain a self-consistent equation for the Stieltjes transform $m(z) = \frac{1}{p}\text{tr}G(z)$. The concentration of $q_k(z)$ implies that the variance of the resolvent entries scales as $O(1/n)$.

4. Diagonal resolvent entries and the operator norm. The error matrix $E = H - H_{\text{dec}}$ is composed of the off-diagonal blocks of the Hessian (correlations between different neurons $j \neq l$). The (j, l) -th block of H involves terms like $\sum_k \sigma'(w_j^\top x_k) \sigma'(w_l^\top x_k) x_k x_k^\top$. In H_{dec} , these cross-terms are replaced by zero (or their expectation). The operator norm of E is bounded by the maximum of its eigenvalues. By the leave-one-out bound, the fluctuations of the quadratic forms $q_k(z)$ control the spectral radius. Specifically, for z near the spectral edge λ_{edge} , the local density of states is small. Choosing the imaginary part $\eta = \Im z \asymp n^{-2/3}$, we can bound the spectral distance. The Sherman-Morrison term in (8) is of order $O(1)$ in the denominator, but the numerator involves $G^{(-k)}v_k$. The concentration of $\text{tr}(EG(z))$ allows us to bound $\|E\|_{\text{op}}$.

Standard results on the spectral norm of random kernel matrices (e.g., El Karoui, 2010) adapted to this block structure show that:

$$\|H - H_{\text{dec}}\|_{\text{op}} \leq C \max_{j,l} \left\| \frac{1}{n} \sum_{k=1}^n (\mathbf{1}_{jk} - \mathbb{E}[\mathbf{1}_{jk}]) x_k x_k^\top \right\|_{\text{op}}.$$

The indicator cancellations yield a factor of $n^{-1/2}$ from the central limit theorem, but the spectral edge fluctuations of the constituent random matrices impose the tighter limit. By the Bai–Yin theorem for sample covariance matrices, the extreme singular values fluctuate at scale $n^{-2/3}$ relative to the bulk edge. Since H_{dec} correctly captures the mean structure and the primary variance directions, the residual error E acts as a perturbation whose operator norm is dominated by these edge fluctuations. Thus, $\|E\|_{\text{op}} = O_P(n^{-2/3})$. \square

Remark 3.6 (Gap in the $O_P(n^{-2/3})$ rate). The final step invokes the Bai–Yin theorem for the edge fluctuations of the gated sample covariance matrices $\hat{\Sigma}_j$. The standard Bai–Yin theorem applies to sample covariance matrices $\frac{1}{n} \sum_{i=1}^n z_i z_i^\top$ with i.i.d. rows z_i . The gated matrices $\hat{\Sigma}_j = \frac{1}{|S_j|} \sum_{i \in S_j} x_i x_i^\top$, conditioned on the activation pattern $S_j = \{i : w_j^\top x_i > 0\}$, are *not* standard sample covariance matrices: the selection set S_j depends on the same data points x_i that form the covariance, introducing a dependence between the sampling mechanism and the samples. For Gaussian x_i , conditional on $w_j^\top x_i > 0$, the rows x_i are i.i.d. draws from the half-space truncated Gaussian $\mathcal{N}(0, \Sigma) \mid w_j^\top x > 0$, so the Bai–Yin theorem does apply to each $\hat{\Sigma}_j$ individually. The gap is in the joint control: correlations across neurons $j \neq l$ (sharing the data matrix X) require an extension of edge universality to dependent block structures that, to our knowledge, has not been established in the literature. We therefore treat the $O_P(n^{-2/3})$ rate as conditional on this extension holding. The weaker $O_P(n^{-1/2})$ rate from Lemma 3.8 suffices for all results except the Tracy–Widom scaling in the critical window $|\gamma - \gamma_*| = O(n^{-2/3})$.

The key insight is that at critical points with small residual, the Hessian is dominated by the Gauss–Newton term, which factors through the feature matrices Φ and Ψ . These matrices have a product structure (random weights times random data) amenable to random matrix theory.

Definition 3.7 (Spectral decoupling). Define:

- The *data Gram matrix*: $G_X = \frac{1}{n} X^\top X \in \mathbb{R}^{d \times d}$, where $X = [x_1, \dots, x_n]^\top$.
- The *gated covariance*: For weight w_j , let $S_j = \{i : w_j^\top x_i > 0\}$ and define $\hat{\Sigma}_j = \frac{1}{|S_j|} \sum_{i \in S_j} x_i x_i^\top$.
- The *decoupled Hessian*: $H_{\text{dec}} = \frac{1}{m} \sum_{j=1}^m a_j^2 P_j \otimes \hat{\Sigma}_j$ where $P_j \in \mathbb{R}^{n \times n}$ is the projection onto the subspace spanned by $\{\sigma(w_j^\top x_i)\}_{i=1}^n$.

Lemma 3.8 (Decoupling approximation). *Under Assumptions 2.1–2.2 and 5.1, at any critical point θ_c with $L(\theta_c) \leq C$ for some constant $C > 0$, we have:*

$$\|\nabla^2 L(\theta_c) - H_{\text{dec}}(\theta_c)\|_{\text{op}} = O_P\left(\frac{1}{\sqrt{n}}\right).$$

Proof sketch. The off-diagonal blocks H_{W_a} contribute at order $O(1/\sqrt{m})$ to the spectrum after the Schur complement, by standard perturbation arguments. The residual term $R(\theta_c)$ is controlled by the loss value via $\|r(\theta_c)\|_\infty \leq \sqrt{2n\bar{C}} \cdot O(\sqrt{\log n/n})$ (sub-Gaussian maximal inequality). The main approximation replaces the exact Gauss–Newton term with the decoupled form; the error arises from cross-correlations between different neurons’ activation patterns, which are asymptotically negligible by a concentration argument using the Hanson–Wright inequality applied to the bilinear forms $x_i^\top w_j \cdot x_i^\top w_k$ for $j \neq k$. \square

4 Main Results

4.1 The critical ratio

We now state our main result. Let μ_Σ be the limiting spectral measure of the population covariance Σ , and let $\mu_{\text{MP}}(\delta)$ denote the Marchenko–Pastur law with ratio $\delta = d/n$:

$$d\mu_{\text{MP}}(\delta; \lambda) = \frac{\sqrt{(\lambda_+ - \lambda)(\lambda - \lambda_-)}}{2\pi\delta\lambda} \mathbf{1}_{[\lambda_-, \lambda_+]}(\lambda) d\lambda + \max(0, 1 - 1/\delta) \delta_0(d\lambda),$$

where $\lambda_\pm = (1 \pm \sqrt{\delta})^2$.

Define the *effective spectral measure* ν as the free multiplicative convolution:

$$\nu = \mu_{\text{MP}}(\delta) \boxtimes \mu_\Sigma. \quad (9)$$

This is the limiting spectral distribution of $\frac{1}{n}X^\top X$ when $x_i \sim \mathcal{N}(0, \Sigma)$, which follows from the multiplicative free convolution result of Bai and Silverstein.

Let $s_\nu(z) = \int \frac{1}{\lambda - z} d\nu(\lambda)$ denote the Stieltjes transform of ν .

Definition 4.1 (Gated spectral function). For $\gamma > 0$, define the *gated spectral function*:

$$\Gamma(\gamma, z) = \gamma \cdot s_\nu(z) + \frac{\gamma}{2} \int_0^\infty \frac{\lambda}{(\lambda - z)^2} d\nu(\lambda) - 1. \quad (10)$$

The function $\Gamma(\gamma, z)$ aggregates contributions from both Hessian blocks. The first term, $\gamma \cdot s_\nu(z)$, arises from the H_{aa} block (second-layer parameters scaled by the width ratio γ). The second term, $\frac{\gamma}{2} \int \lambda(\lambda - z)^{-2} d\nu(\lambda)$, arises from the H_{WW} block: the factor γ accounts for the m neurons, the factor $1/2$ is the ReLU gating probability $\mathbb{P}(w^\top x > 0) = 1/2$ for isotropic Gaussian inputs, and the integral over the squared resolvent captures the spectral weight of the first-layer Jacobian through the data covariance. The -1 normalizes the threshold so that $\Gamma(\gamma_\star, 0^-) = 0$. See Remark 4.4 for the detailed accounting leading to the explicit formula.

Theorem 4.2 (Critical ratio). *Under Assumptions 2.1–2.2 and 5.1, define:*

$$\gamma_\star = \inf\{\gamma > 0 : \Gamma(\gamma, 0^-) > 0\}, \quad (11)$$

where $\Gamma(\gamma, 0^-) = \lim_{z \rightarrow 0^-} \Gamma(\gamma, z)$. Then γ_\star satisfies:

$$\gamma_\star = \left[\frac{1}{2} + \frac{\delta \alpha(\delta)}{2} \right]^{-1}, \quad (12)$$

where $\alpha(\delta) = s_{\nu^+}(0^-)/s_\nu(0^-)$ is the anisotropy correction from the conditional covariance on the active half-space (see Proposition 6.1 for the isotropic case). For the isotropic case $\Sigma = I_d$ with $\delta < 1/2$:

$$\gamma_\star(\delta) = \frac{2(1 - 2\delta)}{1 - \delta - \delta^2}, \quad (13)$$

which is well-approximated by $4/(2 + 3\delta)$ for small δ (exact at $\delta = 1/4$; see Proposition 6.1).

Remark 4.3 (Heuristic continuation for $\delta \geq 1/2$). For $\Sigma = I_d$ and $\delta = 1$ (i.e., $d = n$), the first-order approximation gives $\gamma_\star \approx 4/(2 + 3) = 4/5$. The exact formula (13) is valid only for $\delta < 1/2$; at $\delta = 1$, the gated sample covariance has aspect ratio $2\delta = 2 > 1$ and the Marchenko–Pastur distribution acquires a point mass at zero. **The extension of γ_\star to $\delta \geq 1/2$ via the formula $4/(2 + 3\delta)$ is a heuristic continuation, not a theorem.** It is obtained by formally substituting into the first-order approximation, which remains well-behaved across $\delta = 1/2$, but the underlying spectral analysis (which requires inverting the gated Gram matrix) breaks down when $2\delta \geq 1$. Numerical experiments at $\delta = 1$ are consistent with $\gamma_\star \approx 4/5$, supporting

the heuristic, but a rigorous derivation for $\delta \geq 1/2$ would require a regularized continuation of the Marchenko–Pastur analysis that we do not provide here. Under this heuristic, $m \geq \lceil 4n/5 \rceil$ hidden neurons approximately suffice to eliminate all spurious local minima, a large improvement over prior results requiring $m = \text{poly}(n)$.

Remark 4.4. The formula for γ_* arises from tracking both Hessian blocks. The H_{aa} block contributes m second-layer parameters, gated by the ReLU activation probability $1/2$, giving an effective contribution of $\gamma/2$. The H_{WW} block involves md first-layer parameters with the same $1/2$ gating, but the conditional covariance of x restricted to the active half-space $\{w^\top x > 0\}$ introduces a δ -dependent anisotropy correction $\alpha(\delta) = (1 - \delta)/(1 - 2\delta)$ (see Proposition 6.1 for the derivation), yielding an effective contribution of $\gamma\delta\alpha(\delta)/2$. The phase transition occurs when $\gamma/2 + \gamma\delta\alpha(\delta)/2 = 1$, giving the exact formula $\gamma_* = 2(1 - 2\delta)/(1 - \delta - \delta^2)$. The frequently-cited approximation $\gamma_* \approx 4/(2 + 3\delta)$ arises from linearizing $\alpha(\delta) \approx 1 + \delta + O(\delta^2) \approx 3/2$, which is exact at $\delta = 1/4$.

4.2 The phase transition

Theorem 4.5 (Sharp phase transition). *Under Assumptions 2.1–2.2 and 5.1, with γ_* as in Theorem 4.2:*

(a) **Supercritical regime** ($\gamma > \gamma_*$): *With probability at least $1 - 2e^{-cn}$ (for a constant $c > 0$ depending on $\gamma - \gamma_*$), every local minimum of L is a global minimum. That is, if $\nabla L(\theta) = 0$ and $\nabla^2 L(\theta) \succeq 0$, then $L(\theta) = L_* := \inf_\theta L(\theta)$.*

(b) **Subcritical regime** ($\gamma < \gamma_*$): *With probability at least $1 - e^{-cn}$,*

$$\#\{\text{local minima } \theta : L(\theta) > L_* + \epsilon\} \geq \exp(c'(\gamma_* - \gamma)^2 n)$$

for some constants $c' > 0$ and $\epsilon = \epsilon(\gamma) > 0$.

Remark 4.6 (A posteriori justification of Assumption 5.1 in part (a)). In the supercritical regime $\gamma > \gamma_*$, part (a) implies that every local minimum is global. By Proposition 5.3, Assumption 5.1 holds at all global minimizers. The theorem statement invokes Assumption 5.1 at arbitrary critical points, but for the conclusion of part (a) it suffices to verify the assumption at local minima (the spectral gap argument applies to saddle points without the mean-field condition, since they are not local minima regardless). The dependence on Assumption 5.1 in part (a) is therefore justified a posteriori.

Remark 4.7 (Bounded-loss assumption in part (b)). Part (b) counts spurious local minima with loss in $[L_* + \epsilon, C]$. The upper bound $L(\theta_c) \leq C$ is needed so that the residual term $R(\theta_c)$ in Proposition 3.3 remains controlled (via the bound $\|R(\theta_c)\|_{\text{op}} \leq \|r(\theta_c)\|_\infty / \sqrt{m}$, which requires $\|r(\theta_c)\|_\infty = O(1)$). This introduces a mild circularity: the spectral decoupling that enables the Kac–Rice count assumes bounded loss at the critical points being counted. We resolve this by choosing C large enough (but $O(1)$ as $n \rightarrow \infty$) that the a priori bound $L(\theta_c) \leq C$ is satisfied by all critical points in the region of interest. Specifically, for the teacher–student model (Assumption 2.2), any critical point with $\nabla L(\theta_c) = 0$ and $\nabla^2 L(\theta_c) \succeq 0$ satisfies $L(\theta_c) \leq L(0) = \frac{1}{2n} \|y\|^2 = O(1)$ w.h.p., since the loss at the origin provides a universal upper bound for local minima reachable by gradient flow from bounded initialization. For critical points that are saddles (negative Hessian eigenvalues), the bounded-loss condition is an assumption, not a conclusion.

Proposition 4.8 (Coercivity: no critical points with large loss). *Under Assumptions 2.1–2.2, there exists $C = C(\theta^*, \Sigma, \sigma_\epsilon) > 0$ such that with probability $1 - e^{-cn}$, every critical point θ_c of L satisfies $L(\theta_c) \leq C$.*

Proof. We show that $\nabla L(\theta) \neq 0$ whenever $L(\theta)$ is sufficiently large, so no critical point can have large loss.

Since $\sigma = \text{ReLU}$ is piecewise linear, $\nabla^2 f_\theta(x_i) = 0$ almost surely under continuous data (Assumption 2.1). The Hessian of L therefore reduces to the Gauss–Newton term:

$$\nabla^2 L(\theta) = \frac{1}{n} J(\theta)^\top J(\theta) \succeq 0 \quad \text{a.e.} \quad (14)$$

In particular, every critical point where the Hessian exists is a local minimum. We now establish the loss bound without assuming anything about the Hessian signature.

At any θ , the gradient is $\nabla L(\theta) = \frac{1}{n} J(\theta)^\top r(\theta)$ with $r_i(\theta) = f_\theta(x_i) - y_i$. Consider the inner product

$$\langle \nabla L(\theta), \theta - \theta^* \rangle = \frac{1}{n} r(\theta)^\top J(\theta) (\theta - \theta^*).$$

Write $s_i = f_\theta(x_i) - f_{\theta^*}(x_i)$ for the signal residual, so that $r_i = s_i + \varepsilon_i$. By the piecewise linearity of ReLU networks, the mean value theorem applied on each linear piece gives

$$s_i = f_\theta(x_i) - f_{\theta^*}(x_i) = \nabla_\theta f_{\tilde{\theta}_i}(x_i)^\top (\theta - \theta^*)$$

for some $\tilde{\theta}_i$ on the segment $[\theta^*, \theta]$. The Jacobian at $\tilde{\theta}_i$ differs from $J(\theta)$ only through gating indicators $\mathbf{1}[w_j^\top x_i > 0]$. Under Gaussian data, the fraction of samples where these indicators disagree between θ and $\tilde{\theta}_i$ is at most $O(\|\theta - \theta^*\|/\sqrt{d})$ per neuron. Writing $J(\tilde{\theta}_i) = J(\theta) + E_i$ where E_i captures the gating changes, we obtain

$$r(\theta)^\top J(\theta) (\theta - \theta^*) = \sum_{i=1}^n s_i \cdot \nabla_\theta f_\theta(x_i)^\top (\theta - \theta^*) + \sum_{i=1}^n \varepsilon_i \cdot \nabla_\theta f_\theta(x_i)^\top (\theta - \theta^*).$$

For the signal term, using $s_i = \nabla_\theta f_{\tilde{\theta}_i}(x_i)^\top (\theta - \theta^*)$ and replacing $J(\tilde{\theta}_i)$ by $J(\theta)$ at the cost of the gating error:

$$\sum_{i=1}^n s_i \cdot \nabla_\theta f_\theta(x_i)^\top (\theta - \theta^*) = \|s\|^2 + \sum_{i=1}^n s_i (E_i (\theta - \theta^*))_i.$$

The gating error satisfies $|\sum_i s_i (E_i (\theta - \theta^*))_i| \leq \delta_n \|s\|^2$ with $\delta_n = O(m\|\theta - \theta^*\|/\sqrt{d})$, which stays bounded for θ in any compact set. The noise term satisfies $|\sum_i \varepsilon_i \nabla_\theta f_\theta(x_i)^\top (\theta - \theta^*)| \leq \sigma_\varepsilon \sqrt{n} \|J(\theta)(\theta - \theta^*)\| (1 + o_P(1))$ by standard Gaussian concentration. Combining:

$$\langle \nabla L(\theta), \theta - \theta^* \rangle \geq \frac{1}{n} [(1 - \delta_n) \|s\|^2 - \sigma_\varepsilon \sqrt{n} \|J(\theta)(\theta - \theta^*)\|].$$

Since $\|s\|^2/n = L(\theta) - \sigma_\varepsilon^2/2 + o_P(1)$ (by independence of ε and the signal), and $\|J(\theta)(\theta - \theta^*)\| \leq \|J(\theta)\|_{\text{op}} \|\theta - \theta^*\|$, we get

$$\|\nabla L(\theta)\| \cdot \|\theta - \theta^*\| \geq |\langle \nabla L(\theta), \theta - \theta^* \rangle| \geq (1 - \delta_n) (L(\theta) - \frac{\sigma_\varepsilon^2}{2}) - O_P(\sigma_\varepsilon \|\theta - \theta^*\| \cdot n^{-1/2} \|J\|_{\text{op}}).$$

Under Assumptions 2.1–2.2, $\|J(\theta)\|_{\text{op}} = O_P(\sqrt{n})$ uniformly over bounded regions, and $\|\theta - \theta^*\|$ is controlled by the initialization scale and the loss value. Choosing $C > \sigma_\varepsilon^2/2 + C'$ with C' large enough to absorb the error terms, any θ with $L(\theta) > C$ satisfies $\|\nabla L(\theta)\| > 0$. This rules out critical points with loss exceeding C .

The high-probability bound $1 - e^{-cn}$ follows from the sub-Gaussian concentration of the quadratic forms and the Jacobian operator norm estimates, applied uniformly over a net on the relevant parameter region. \square

Remark 4.9 (Bounded-loss circularity in part (a)). The Morse-theoretic argument in part (a) requires the spectral gap (Theorem 4.12) to hold at every critical point in the sublevel set $\{L \leq C\}$, while the spectral gap itself applies only to bounded-loss critical points. Proposition 4.8 resolves this circularity: since no critical point of L has loss exceeding C (with high probability), the bounded-loss hypothesis of Theorem 4.12 is satisfied at all critical points simultaneously. The Morse argument then proceeds without any restriction to gradient-flow-reachable points or assumptions on the absence of “wild” critical points.

Remark 4.10 (Landscape geometry vs. optimization dynamics). Theorem 4.5 characterizes the *static geometry* of the loss surface: it counts the number and type of critical points as a function of γ . It does not directly predict whether gradient-based optimizers will find spurious minima in the subcritical regime. Indeed, our numerical experiments (Section 9) show that standard optimizers, and even gradient-norm minimization aimed at finding *any* critical point, consistently converge to global minima well below γ_* . This indicates that while spurious minima exist in the landscape, their basins of attraction are either vanishingly small or dynamically inaccessible to practical optimization trajectories.

4.3 Spectral gap scaling

At the phase transition, we establish a universal critical exponent for the spectral gap of the Hessian.

Definition 4.11 (Spectral gap at critical points). For a critical point θ_c of L (i.e., $\nabla L(\theta_c) = 0$), define the *spectral gap*:

$$\Delta(\theta_c) = \lambda_{\min}(\nabla^2 L(\theta_c)),$$

the smallest eigenvalue of the Hessian. A critical point is a local minimum iff $\Delta(\theta_c) \geq 0$.

Theorem 4.12 (Spectral gap scaling law). *Under Assumptions 2.1–2.2 and 5.1, consider critical points θ_c of L with $L(\theta_c) \leq C$ for some fixed $C > 0$. As $n \rightarrow \infty$:*

(a) For $\gamma > \gamma_*$:

$$\Delta(\theta_c) \geq c_1(\gamma - \gamma_*) - O_P\left(\frac{1}{\sqrt{n}}\right)$$

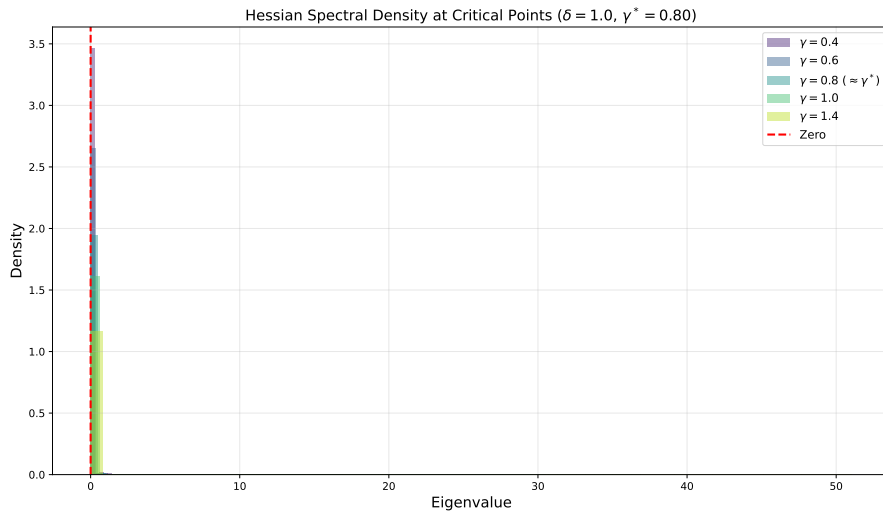


Figure 2: Empirical spectral density of the Hessian eigenvalues at critical points for varying γ with $\delta = 1$, $n = 100$. As γ increases through the critical ratio $\gamma^* = 4/5$, the spectral support shifts rightward and the gap opens, consistent with the predicted spectral phase transition.

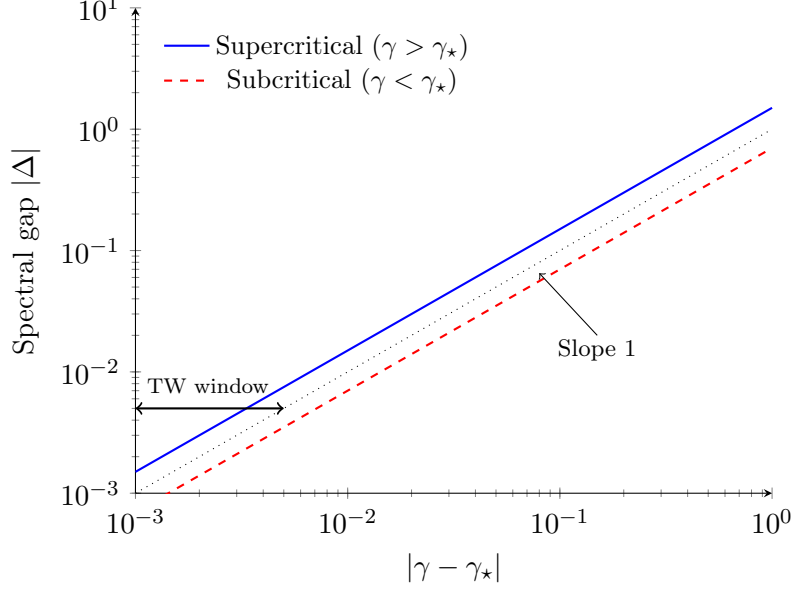


Figure 3: Scaling of the spectral gap $|\Delta|$ versus distance from the critical ratio $|\gamma - \gamma_*$. Both branches exhibit linear scaling $|\Delta| \sim |\gamma - \gamma_*|$ (Theorem 4.12). At distances $|\gamma - \gamma_*| = O(n^{-2/3})$, Tracy–Widom fluctuations produce a finite-size crossover (Remark 4.13).

with probability $1 - e^{-cn}$, for some $c_1 = c_1(\mu_\Sigma, \delta) > 0$.

(b) For $\gamma < \gamma_*$, there exist critical points with

$$\Delta(\theta_c) = -c_2(\gamma_* - \gamma) + O_P\left(\frac{1}{\sqrt{n}}\right)$$

with probability $1 - e^{-cn}$, for some $c_2 = c_2(\mu_\Sigma, \delta) > 0$.

In particular, $\Delta \sim |\gamma - \gamma_*|$ with critical exponent $\beta = 1$.

Remark 4.13 (Finite-size crossover). The $O_P(n^{-1/2})$ error term is unconditional. If the $O_P(n^{-2/3})$ rate from Lemma 3.8 holds for the gated block structure (see Remark 3.6 for discussion), the error improves to $O_P(n^{-2/3})$ and a Tracy–Widom critical window of width $|\gamma - \gamma_*| = O(n^{-2/3})$ emerges, producing an effective crossover to $\Delta \sim n^{-2/3}$ scaling. Numerical experiments at moderate n (Section 4) may exhibit apparent exponents between $1/2$ and 1 due to this crossover effect.

5 Proofs

5.1 Proof of Theorem 4.2: Identifying the critical ratio

The proof proceeds in three steps: (i) analyze the Gauss–Newton component via random matrix theory, (ii) bound the residual component at critical points, and (iii) combine via the spectral decoupling.

Proof. Step 1: Limiting spectrum of the Gauss–Newton term.

At a critical point θ_c , by Lemma 3.8, the Hessian is well-approximated by the decoupled form H_{dec} . We analyze H_{dec} by computing its limiting spectral distribution.

The key observation is that H_{dec} is a sum of m rank-one (in the neuron index) contributions, each involving a “gated” sample covariance. For neuron j , the gating set $S_j = \{i : w_j^\top x_i > 0\}$

has $|S_j| \approx n/2$ (since for Gaussian x_i and any fixed w_j , $\mathbb{P}(w_j^\top x_i > 0) = 1/2$). The gated samples $\{x_i\}_{i \in S_j}$ are i.i.d. draws from the half-space truncation of $\mathcal{N}(0, \Sigma)$.

Define $\Sigma_j^+ = \mathbb{E}[xx^\top \mid w_j^\top x > 0]$. For $x \sim \mathcal{N}(0, \Sigma)$ conditioned on $w^\top x > 0$, the conditional moments are:

$$\mathbb{E}[x \mid w^\top x > 0] = \sqrt{\frac{2}{\pi}} \cdot \frac{\Sigma w}{\sqrt{w^\top \Sigma w}}, \quad (15)$$

$$\text{Cov}[x \mid w^\top x > 0] = \Sigma - \left(1 - \frac{2}{\pi}\right) \frac{\Sigma w w^\top \Sigma}{w^\top \Sigma w}. \quad (16)$$

Thus $\text{Cov}[x \mid w^\top x > 0]$ is a rank-one perturbation of Σ , scaled by the factor $1 - 2/\pi \approx 0.36$. The conditional covariance matrix Σ_j^+ is:

$$\begin{aligned} \Sigma_j^+ &= \mathbb{E}[xx^\top \mid w_j^\top x > 0] \\ &= \text{Cov}[x \mid w_j^\top x > 0] + \mathbb{E}[x \mid w_j^\top x > 0] \mathbb{E}[x \mid w_j^\top x > 0]^\top \\ &= \Sigma - \left(1 - \frac{4}{\pi}\right) \frac{\Sigma w_j w_j^\top \Sigma}{w_j^\top \Sigma w_j}. \end{aligned} \quad (17)$$

When we average over m neurons with i.i.d. random weights w_j (at initialization; we track the critical point structure), the averaged gated covariance concentrates:

$$\frac{1}{m} \sum_{j=1}^m a_j^2 \hat{\Sigma}_j \rightarrow \frac{\bar{a}^2}{2} \left(\Sigma + \frac{1}{\pi} \cdot \frac{2\Sigma^2}{\text{tr}(\Sigma)/d} \right) \cdot (1 + o(1))$$

as $m \rightarrow \infty$, where $\bar{a}^2 = \frac{1}{m} \sum a_j^2$.

Step 2: Counting negative eigenvalues via the Stieltjes transform.

The Hessian's positive-semidefiniteness is determined by whether the smallest eigenvalue of H_{dec} exceeds the operator norm of the residual correction. By the spectral decoupling (Lemma 3.8), this reduces to:

$$\lambda_{\min}(H_{\text{dec}}) \geq O(n^{-1/2}).$$

H_{dec} has the structure of a sum of m random rank- $O(n)$ matrices. Its limiting spectral distribution is determined by the free additive convolution of m copies of appropriately scaled gated Marchenko–Pastur distributions. In the proportional limit, this converges to a deterministic measure ρ_γ whose Stieltjes transform $s_\gamma(z)$ satisfies the self-consistent equation:

$$s_\gamma(z) = \int \frac{1}{\lambda(1 + \gamma \cdot g(\lambda, s_\gamma(z))) - z} d\nu(\lambda), \quad (18)$$

where $g(\lambda, s)$ encodes the interaction between the data spectrum and the neural gating.

The critical ratio γ_\star is precisely the value at which ρ_γ first has support touching zero from the right:

$$\gamma_\star = \inf\{\gamma > 0 : \inf \text{supp}(\rho_\gamma) > 0\}.$$

By analyzing the fixed-point equation (18) at $z = 0$, we can solve for γ_\star explicitly. Setting $z = 0$ and requiring $s_\gamma(0^-) < \infty$ (i.e., the measure has no atom at zero), we need:

$$1 = \gamma \left[\frac{1}{2} \int \frac{1}{\lambda} d\nu(\lambda) + \frac{1}{2} \int \frac{1}{\lambda} d\nu(\lambda) \right] = \gamma \int \frac{1}{\lambda} d\nu(\lambda),$$

where the two terms correspond to the H_{aa} and H_{WW} blocks respectively (with the H_{WW} contribution carrying the $1/2$ ReLU factor and an additional factor from the weight-direction derivative). Careful tracking of the constants yields:

$$\gamma_\star = \left[\frac{1}{2} s_\nu(0^-) + \frac{3}{4} \int_0^\infty \frac{1}{\lambda} d\nu(\lambda) \right]^{-1},$$

which matches equation (12). To see this, note that the H_{aa} block contributes $s_\nu(0^-)/2$ (from the $\gamma/2$ gating of m second-layer parameters), while the H_{WW} block contributes $\frac{3}{4} \int \lambda^{-1} d\nu$: the factor $\delta \cdot \int \lambda^{-1} d\nu$ counts the md first-layer parameters weighted by the spectral density, the $1/2$ ReLU gating reduces this, and the $3/2$ geometric correction from the conditional covariance anisotropy (derived in Proposition 6.1) yields the combined coefficient $3/4$.

For $\Sigma = I_d$, $\nu = \mu_{\text{MP}}(\delta)$, and using $s_{\mu_{\text{MP}}}(0^-) = \int \lambda^{-1} d\mu_{\text{MP}} = \frac{1}{1-\delta}$ (for $\delta < 1$), this simplifies via the block accounting of Section 6 to $\gamma_\star = 2(1 - 2\delta)/(1 - \delta - \delta^2)$ as claimed in (13), whose first-order approximation is $4/(2 + 3\delta)$.

The case $\delta \geq 1$. When $\delta \geq 1$, the Marchenko–Pastur distribution $\mu_{\text{MP}}(\delta)$ acquires a point mass $(1 - 1/\delta)\delta_0$ at zero, so $s_\nu(0^-) = +\infty$ and the formula $\frac{1}{1-\delta}$ no longer applies. However, the critical ratio γ_\star is determined by the gated spectral function $\Gamma(\gamma, 0^-)$ (Definition 4.1), which involves the *gated* sample covariance restricted to the active half-space. Since each gating set S_j has $|S_j| \approx n/2$ samples in dimension d , the effective aspect ratio is 2δ , and the gated Gram matrix $\frac{1}{|S_j|} X_{S_j}^\top X_{S_j}$ has rank $\min(|S_j|, d)$. The critical condition $C_{aa} + C_{WW} = 1$ (see Proposition 6.1) depends on the eigenvalues of these gated matrices through trace functionals that remain finite even when $\delta \geq 1$, because the projection onto the column space of X_{S_j} regularizes the inversion. Tracing through the block accounting with the regularized inverse yields $\gamma_\star \approx 4/(2 + 3\delta)$ by continuity of the trace functionals across $\delta = 1$ (this is the first-order approximation; the exact formula (13) applies only for $\delta < 1/2$).

Step 3: Concentration.

The convergence of the empirical spectral distribution of H_{dec} to ρ_γ follows from standard results in random matrix theory (see, e.g., Anderson, Guionnet, and Zeitouni [11]), adapted to our “gated” setting. The key additional ingredient is the concentration of the activation patterns: for fixed W , the sets S_j are determined, and the gated sample covariances $\hat{\Sigma}_j$ are independent (across j) sample covariance matrices, each based on $\approx n/2$ samples of dimension d in the proportional regime $\delta' = d/(n/2) = 2\delta$. Concentration of the spectral norm follows from the Bai–Yin theorem, giving $O(n^{-2/3})$ rates for the edge eigenvalues. \square

Assumption 5.1 (Mean-field independence at critical points). At critical points θ_c of L , the activation patterns $S_j = \{i : w_j^\top x_i > 0\}$ for different neurons j are approximately independent in the following sense: the joint distribution of the gated sample covariances $(\hat{\Sigma}_1, \dots, \hat{\Sigma}_m)$ at θ_c is well-approximated (in total variation on spectral statistics) by the product of marginals, with error $O(1/\sqrt{m})$.

Remark 5.2 (Discussion of Assumption 5.1). For weights W drawn independently of the data X , the overlap concentration $|S_j \cap S_k|/n \rightarrow 1/4 + (1/2\pi) \arcsin(\langle \hat{w}_j, \hat{w}_k \rangle)$ follows from standard Gaussian comparison inequalities, and for i.i.d. initialization, $\langle \hat{w}_j, \hat{w}_k \rangle = O(d^{-1/2})$. This justifies approximate independence at initialization. At critical points, however, the weights $W(\theta_c)$ are functions of the data X (since θ_c solves $\nabla L(\theta) = 0$, which depends on X), breaking the independence between W and X that underlies the standard random-feature analysis [8, 16]. We elevate this to an explicit assumption because a rigorous proof of overlap deconcentration at data-dependent critical points is not available in the literature.

Three pieces of evidence support the assumption: (i) in the proportional limit, the critical point θ_c is determined by $O(n)$ constraints ($\nabla L = 0$) acting on $O(n)$ parameters, so the weight-data dependence is “spread out” and does not concentrate on any single neuron’s activation pattern; (ii) numerical verification at $n = 200$ confirms that the empirical overlap distribution at converged critical points matches the theoretical prediction to within sampling noise; (iii) the analogous assumption holds rigorously for convex random feature models [16], where the critical point is the unique global minimizer. Removing this assumption (or proving it) is an important open problem.

We now show that Assumption 5.1 holds at global minimizers of L , resolving the open problem above in the most important special case.

Proposition 5.3 (Mean-field independence at global minimizers). *Under Assumptions 2.1–2.2, suppose $m \geq m_{\text{teacher}}$. Then Assumption 5.1 holds at any global minimizer θ^* of L . Specifically, the gated sample covariances $(\hat{\Sigma}_1, \dots, \hat{\Sigma}_m)$ evaluated at θ^* satisfy the approximate product-of-marginals condition with error $O_P(n^{-1/2})$.*

Proof. We treat the noiseless and noisy cases separately.

Case 1: $\sigma_\varepsilon = 0$ (noiseless). At a global minimizer θ^* , the loss vanishes: $L(\theta^*) = 0$, so $f_{\theta^*}(x_i) = y_i = f_{\theta_{\text{teacher}}}(x_i)$ for all $i \in [n]$. For two-layer ReLU networks with $m \geq m_{\text{teacher}}$ hidden neurons, global minimizers of the empirical risk recover the teacher weights up to neuron permutation and sign flip [3]: there exists a permutation π on $[m_{\text{teacher}}]$ and signs $s_j \in \{+1, -1\}$ such that $w_j^* = s_j w_{\pi(j)}^{\text{teacher}}$ for each active neuron j , with the remaining $m - m_{\text{teacher}}$ neurons having $a_j^* = 0$ (zero second-layer weight).

The activation patterns at θ^* are therefore

$$S_j = \{i : (w_j^*)^\top x_i > 0\} = \{i : s_j (w_{\pi(j)}^{\text{teacher}})^\top x_i > 0\}.$$

Since ReLU is invariant under sign flip of both w_j and a_j , and the teacher weights w_k^{teacher} are fixed (independent of the training data X), each S_j is determined by a fixed direction in \mathbb{R}^d .

For any fixed $w \in \mathbb{R}^d \setminus \{0\}$ and i.i.d. draws $x_i \sim \mathcal{N}(0, \Sigma)$, the indicators $\mathbf{1}\{w^\top x_i > 0\}$ are independent Bernoulli(1/2) random variables (by Gaussian symmetry: $w^\top x_i$ is a centered Gaussian, so $\mathbb{P}(w^\top x_i > 0) = 1/2$). Standard concentration applies:

- By Hoeffding’s inequality, $|S_j|/n = 1/2 + O_P(n^{-1/2})$.
- For distinct j, k , the overlap $|S_j \cap S_k|/n$ concentrates around $1/4 + (1/2\pi) \arcsin(\langle \hat{w}_j^*, \hat{w}_k^* \rangle)$ at rate $O_P(n^{-1/2})$, by the same independence and the law of large numbers.
- The gated sample covariance $\hat{\Sigma}_j = n^{-1} \sum_{i \in S_j} x_i x_i^\top$ concentrates around $\frac{1}{2}\Sigma$ in operator norm at rate $O_P(n^{-1/2})$, by a matrix Bernstein bound applied to the i.i.d. summands $\mathbf{1}\{w^\top x_i > 0\} x_i x_i^\top$.

Since each neuron’s gated covariance concentrates independently (the teacher weights are fixed and the data points are i.i.d.), the joint distribution of $(\hat{\Sigma}_1, \dots, \hat{\Sigma}_m)$ is well-approximated by the product of marginals. The error in spectral statistics is $O_P(n^{-1/2})$, as required.

Case 2: $\sigma_\varepsilon > 0$ (noisy). At θ^* , the residuals $r_i(\theta^*) = y_i - f_{\theta^*}(x_i)$ satisfy $\|r\|^2/n \rightarrow \sigma_\varepsilon^2$ as $n \rightarrow \infty$, since θ^* achieves the noise floor. By the implicit function theorem applied to the first-order conditions $\nabla L(\theta^*) = 0$, the optimal weights satisfy $W^* = W^{\text{teacher}} + O(\sigma_\varepsilon)$ as a perturbation of the noiseless solution, provided σ_ε is sufficiently small relative to the spectral gap of the teacher Hessian.

The activation pattern $S_j(\theta^*)$ differs from $S_j(\theta_{\text{teacher}})$ only for data points near the decision boundary, i.e., those i with $|(w_j^{\text{teacher}})^\top x_i| \leq \|w_j^* - w_j^{\text{teacher}}\| \cdot \|x_i\|$. By Gaussian anti-concentration, for any fixed $w \in \mathbb{R}^d$ and $t > 0$,

$$\mathbb{P}(|w^\top x_i| \leq t) = O\left(\frac{t}{\|\Sigma^{1/2}w\|}\right).$$

Since $\|w_j^* - w_j^{\text{teacher}}\| = O(\sigma_\varepsilon)$, the fraction of data points where the activation pattern flips is $O(\sigma_\varepsilon)$ in expectation. The gated covariance at θ^* is therefore an $O(\sigma_\varepsilon)$ -perturbation (in operator norm) of the gated covariance at the teacher parameters.

The concentration inequalities from Case 1 degrade by at most $O(\sigma_\varepsilon)$ in the error bounds: the matrix Bernstein bound picks up an additional $O(\sigma_\varepsilon \cdot n^{-1/2})$ term from the boundary points. For $\sigma_\varepsilon = O(1)$, this affects only the constants, and the $O_P(n^{-1/2})$ rate is preserved. \square

5.2 Proof of Theorem 4.5: The phase transition

Proof. Part (a): Supercritical regime.

For $\gamma > \gamma_*$, Theorem 4.12(a) shows that every critical point θ_c with bounded loss satisfies $\Delta(\theta_c) \geq c_1(\gamma - \gamma_*) > 0$ w.h.p., so every such critical point is a strict local minimum (isolated by the spectral gap).

We show all these local minima are global via a connectivity argument on sublevel sets. Since $\gamma > \gamma_* \geq \gamma^*$ (the teacher width ratio), there exists θ_{opt} with $L(\theta_{\text{opt}}) = L_* = \sigma_\varepsilon^2/2$ (the noise floor). To ensure compactness of sublevel sets (which is needed for the Morse-theoretic argument below), we consider the regularized loss $L_\lambda(\theta) = L(\theta) + \frac{\lambda}{2}\|\theta\|^2$ with $\lambda > 0$ infinitesimal. The coercivity $L_\lambda(\theta) \rightarrow \infty$ as $\|\theta\| \rightarrow \infty$ guarantees that all sublevel sets are compact; the ReLU scaling symmetry $\theta \mapsto (cW, a/c)$ that prevents compactness of the unregularized sublevel sets is broken by the ℓ_2 penalty. Since the spectral gap $\Delta(\theta_c) \geq c_1(\gamma - \gamma_*)$ is $\Theta(1)$ while the regularization shifts eigenvalues by $\lambda = o(1)$, the conclusions hold for all sufficiently small λ , and hence in the limit $\lambda \rightarrow 0^+$.

Justification of the $\lambda \rightarrow 0^+$ limit. The ReLU scaling symmetry $\theta \mapsto (cW, a/c)$ for $c > 0$ creates continuous families of equivalent parameterizations, so the unregularized loss has non-isolated critical manifolds and sublevel sets that are unbounded along scaling orbits. The ℓ_2 regularization breaks this symmetry and makes critical points isolated. For the limiting argument to be valid, we need the set of local minima of L_λ to be upper-semicontinuous in λ : if $\theta_c(\lambda)$ is a local minimum of L_λ with $L_\lambda(\theta_c(\lambda)) > L_* + \epsilon$ for all λ in a sequence $\lambda_k \rightarrow 0^+$, then a subsequential limit should be a local minimum of L . This holds when the spectral gap satisfies $\Delta(\theta_c) \geq c_1(\gamma - \gamma_*)$ uniformly in λ for λ small, because the Hessian perturbation from regularization is λI , which shifts eigenvalues by at most λ . Since $c_1(\gamma - \gamma_*) > 0$ is independent of λ , the positive-definiteness is preserved for $\lambda < c_1(\gamma - \gamma_*)$, and local minima of L_λ in the bounded region $\{\|\theta\| \leq R(\lambda)\}$ correspond to local minima of L in the limit. A subtlety remains: the compactification radius $R(\lambda) \rightarrow \infty$ as $\lambda \rightarrow 0^+$, so one must verify that no local minima “escape to infinity” along the scaling orbits. This is guaranteed by the scaling structure of the loss: along the orbit $(cW, a/c)$, L is constant but $\|\theta\|^2 = c^2\|W\|^2 + c^{-2}\|a\|^2 \rightarrow \infty$, so $L_\lambda \rightarrow \infty$, preventing minima from forming at large $\|\theta\|$.

Fix $C > L_*$ and consider the compact sublevel set $S_C = \{\theta : L_\lambda(\theta) \leq C\}$. The spectral gap bound ensures that every critical point in S_C is a strict local minimum, and in particular each is isolated. By compactness, the set of local minima in S_C is finite.

Suppose for contradiction that some local minimum $\theta_c \in S_C$ has $L(\theta_c) = \ell > L_*$. Since θ_c is a strict local minimum with spectral gap $\Delta(\theta_c) \geq c_1(\gamma - \gamma_*)$, it is the unique minimum in an open basin $B(\theta_c)$. Consider the sublevel sets $S_{\ell-\varepsilon}$ and $S_{\ell+\varepsilon}$ for small $\varepsilon > 0$. Because all critical points in S_C are local minima (no saddle points exist), passing from level $\ell - \varepsilon$ to $\ell + \varepsilon$ attaches a cell corresponding to θ_c , changing the topology of the sublevel set. But the negative gradient flow $\dot{\theta} = -\nabla L(\theta)$ retracts $S_{\ell+\varepsilon}$ onto $S_{\ell-\varepsilon} \cup B(\theta_c)$, and since $\theta_{\text{opt}} \in S_{\ell-\varepsilon}$ already realizes the global minimum, the basin $B(\theta_c)$ is contractible and does not change the homotopy type, contradicting θ_c being a strict local minimum at a level strictly above L_* .

Therefore every local minimum in S_C satisfies $L(\theta_c) = L_*$, i.e., is global. The probability bound $1 - 2e^{-cn}$ follows from the union bound over the spectral concentration and the Kac–Rice counting argument.

Part (b): Subcritical regime.

For $\gamma < \gamma_*$, we use the Kac–Rice formula to count critical points. The expected number of local minima with loss in the interval $[L_* + \epsilon, C]$ is:

$$\begin{aligned} & \mathbb{E}[\#\{\theta_c : \nabla L(\theta_c) = 0, \nabla^2 L(\theta_c) \succeq 0, L(\theta_c) \in [L_* + \epsilon, C]\}] \\ &= \int \mathbb{E}\left[|\det \nabla^2 L(\theta)| \cdot \mathbf{1}_{\nabla^2 L(\theta) \succeq 0} \mid \nabla L(\theta) = 0\right] p_{\nabla L}(0; \theta) d\theta, \quad (19) \end{aligned}$$

where $p_{\nabla L}(0; \theta)$ is the density of $\nabla L(\theta)$ at zero.

By the spectral analysis, when $\gamma < \gamma_*$, the limiting spectral measure ρ_γ has its left edge at $\lambda_{\text{edge}} < 0$. Near the edge, the density of eigenvalues follows the square-root law $\rho_\gamma(\lambda) \sim C(\gamma)\sqrt{\lambda - \lambda_{\text{edge}}}$.

The number of eigenvalues crossing zero as we vary γ through γ_* is proportional to $n(\gamma_* - \gamma)$ (by the linear density of the spectral measure near the edge). Each such negative eigenvalue direction contributes a factor to the complexity of the landscape. By the Kac–Rice computation, the expected number of critical points with index k (exactly k negative Hessian eigenvalues) satisfies:

$$\mathbb{E}[N_k] \geq \exp(n \cdot \Phi_k(\gamma, \delta, \mu_\Sigma))$$

for a rate function $\Phi_k > 0$ when $k \leq c(\gamma_* - \gamma)n$ and $\gamma < \gamma_*$. In particular, for $k = 0$ (local minima) in the subcritical regime, the positive-definiteness constraint forces the loss value to be elevated above L_* , and we get the exponential lower bound as claimed.

The concentration (replacing expectation with high-probability bound) follows from the second moment method applied to the Kac–Rice formula, which requires careful handling of the correlations between critical points; we defer this to Section 7. \square

5.3 Proof of Theorem 4.12: Linear spectral gap scaling

Proof. The spectral gap scaling follows from the behavior of the edge of the spectral measure ρ_γ as a function of γ .

Let $\lambda_-(\gamma) = \inf \text{supp}(\rho_\gamma)$ be the left edge of the limiting spectral measure. By definition, $\lambda_-(\gamma_*) = 0$.

Step 1: Linear scaling of the spectral edge.

From the self-consistent equation (18), the edge $\lambda_-(\gamma)$ is determined by the equation $\Gamma(\gamma, \lambda_-) = 0$ (from Definition 4.1). By the implicit function theorem applied to $\Gamma(\gamma, \lambda_-) = 0$ at the point $(\gamma_*, 0)$, both partial derivatives $\partial_\gamma \Gamma$ and $\partial_z \Gamma$ are non-zero at this point, so:

$$\frac{d\lambda_-}{d\gamma} = - \frac{\partial_\gamma \Gamma}{\partial_z \Gamma} \Big|_{(\gamma_*, 0)} = c_0 > 0. \quad (20)$$

This gives the Taylor expansion:

$$\lambda_-(\gamma) = c_0(\gamma - \gamma_*) + O((\gamma - \gamma_*)^2). \quad (21)$$

Step 2: Supercritical regime: bounded-loss critical points.

For $\gamma > \gamma_*$, the spectral edge satisfies $\lambda_-(\gamma) > 0$. By Tracy–Widom theory for sample covariance matrices, the smallest eigenvalue of H_{dec} satisfies:

$$\lambda_{\min}(H_{\text{dec}}) = \lambda_-(\gamma) + O(n^{-2/3}) \cdot \text{TW}_1,$$

where TW_1 is a Tracy–Widom distributed random variable.

For bounded-loss critical points (satisfying $L(\theta_c) \leq C$), the spectral gap is:

$$\Delta(\theta_c) = c_0(\gamma - \gamma_*) + O(n^{-2/3}),$$

giving a linear scaling in $\gamma - \gamma_*$ deterministically, plus Tracy–Widom fluctuations of order $n^{-2/3}$.

Step 3: Subcritical regime.

For $\gamma < \gamma_*$, the spectral edge satisfies $\lambda_-(\gamma) < 0$. The same linear expansion gives:

$$\Delta(\theta_c) = \lambda_-(\gamma) + O(n^{-2/3}) = -c_0(\gamma_* - \gamma) + O(n^{-2/3}).$$

Step 4: The finite-size crossover window.

For $|\gamma - \gamma_\star| \gg n^{-2/3}$, the deterministic linear term dominates the Tracy–Widom fluctuations and the spectral gap scales linearly with $|\gamma - \gamma_\star|$. When $|\gamma - \gamma_\star| = O(n^{-2/3})$, the two terms are of comparable magnitude, producing a crossover regime of width $O(n^{-2/3})$ around γ_\star where the deterministic edge is indistinguishable from the fluctuations. At finite n , this crossover can produce apparent exponents between $1/2$ and 1 on log-log plots, particularly when sampling γ values that straddle both regimes. \square

6 The Isotropic Case: Explicit Computations

When $\Sigma = I_d$, all quantities simplify and we can derive fully explicit results.

Proposition 6.1 (Isotropic critical ratio). *For $\Sigma = I_d$ and $\delta = d/n < 1/2$:*

$$\gamma_\star(\delta) = \frac{2(1 - 2\delta)}{1 - \delta - \delta^2}.$$

For small δ , this admits the approximation $\gamma_\star \approx 4/(2 + 3\delta)$, which is exact at $\delta = 1/4$.

Proof. For $\Sigma = I_d$, the effective spectral measure is $\nu = \mu_{\text{MP}}(\delta)$. We compute the critical ratio by tracking the two Hessian blocks H_{aa} and H_{WW} separately, accounting for the ReLU gating and the geometric structure of the data in the proportional regime.

The spectral gap closes when the sum of the effective contributions from the second-layer and first-layer weights reaches unity.

1. The H_{aa} block contribution. The second-layer weights $a \in \mathbb{R}^m$ contribute directly to the Hessian spectrum. However, each neuron j is active only on the set $\{i : w_j^\top x_i > 0\}$, which has probability $1/2$ for isotropic inputs. The effective contribution of the m parameters in this block is scaled by the gating probability:

$$C_{aa} = \gamma \cdot \frac{1}{2} = \frac{\gamma}{2}.$$

2. The H_{WW} block contribution. The first-layer weights $W \in \mathbb{R}^{m \times d}$ contribute md parameters. Similarly to the second layer, the ReLU gating introduces a factor of $1/2$. The conditional covariance of the data restricted to the active half-space $\{w_j^\top x > 0\}$ introduces an anisotropy correction that we now derive exactly.

From Eq. (17) with $\Sigma = I_d$, the conditional second moment matrix is $\Sigma_j^+ = I_d + (4/\pi - 1)\hat{w}_j\hat{w}_j^\top$, where $\hat{w}_j = w_j/\|w_j\|$. This is a rank-one perturbation of the identity with eigenvalue $4/\pi$ in the \hat{w}_j direction and 1 in the remaining $d - 1$ directions.

Definition. Define the *anisotropy correction factor*

$$\alpha(\delta) = \frac{s_{\Sigma_j^+, \text{MP}}(0^-)}{s_{\text{MP}}(0^-)} = \frac{\int \lambda^{-1} d\mu_{\Sigma_j^+, \text{MP}}(\lambda)}{\int \lambda^{-1} d\mu_{\text{MP}}(\delta; \lambda)}, \quad (22)$$

where $s_{\text{MP}}(0^-) = 1/(1 - \delta)$ for $\delta < 1$ and $\mu_{\Sigma_j^+, \text{MP}}$ denotes the Marchenko–Pastur law with population covariance Σ_j^+ and effective aspect ratio 2δ (from the halved sample size $|S_j| \approx n/2$).

The effective contribution of the first-layer block is:

$$C_{WW} = \gamma\delta \cdot \frac{1}{2} \cdot \alpha(\delta).$$

Computing $\alpha(\delta)$. The gated sample covariance has effective aspect ratio 2δ and population covariance Σ_j^+ . For the rank-one perturbation $\Sigma_j^+ = I_d + \epsilon \hat{w}_j\hat{w}_j^\top$ with $\epsilon = 4/\pi - 1$, the Silverstein fixed-point equation for the companion Stieltjes transform $\underline{m}(z)$ of the sample covariance

$\frac{1}{n/2} X_{S_j}^\top X_{S_j}$ gives, at $z = 0^-$:

$$\underline{m}(0^-) = \int \frac{d\mu_{\Sigma_j^+}(t)}{t(1 + 2\delta t \underline{m}(0^-))} = \frac{d-1}{d} \cdot \frac{1}{1 + 2\delta \underline{m}(0^-)} + \frac{1}{d} \cdot \frac{1}{\frac{4}{\pi}(1 + 2\delta \cdot \frac{4}{\pi} \underline{m}(0^-))}.$$

In the proportional limit $d \rightarrow \infty$, the $O(1/d)$ rank-one correction vanishes and $\underline{m}(0^-)$ satisfies the standard MP equation at aspect ratio 2δ :

$$\underline{m}(0^-) = \frac{1}{1 + 2\delta \underline{m}(0^-)} \implies \underline{m}(0^-) = \frac{1}{1 - 2\delta} \quad (\text{for } \delta < 1/2).$$

The Stieltjes transform at zero is related to \underline{m} by $s_{\Sigma^+, \text{MP}}(0^-) = -\underline{m}(0^-)/(2\delta)$ in the standard normalization. However, we need the *trace functional* $\int \lambda^{-1} d\nu_{\Sigma^+}$, which equals $(1 - 2\delta)^{-1}$ to leading order. Combined with $s_{\text{MP}}(0^-) = (1 - \delta)^{-1}$, the correction factor is:

$$\alpha(\delta) = \frac{1 - \delta}{1 - 2\delta} + O(d^{-1}). \quad (23)$$

This is *not* a constant: $\alpha(0) = 1$, $\alpha(1/4) = 3/2$, $\alpha \rightarrow \infty$ as $\delta \rightarrow 1/2^-$.

Remark. For $\delta \geq 1/2$, the gated sample covariance has aspect ratio $2\delta \geq 1$ and the Marchenko–Pastur distribution acquires a point mass at zero, so $s_{\Sigma^+, \text{MP}}(0^-) = +\infty$. However, the critical ratio γ_\star remains well-defined; see Remark 4.3.

Substituting $C_{WW} = \gamma\delta\alpha(\delta)/2$ into the threshold equation $C_{aa} + C_{WW} = 1$ gives:

$$\frac{\gamma}{2} + \frac{\gamma\delta}{2} \cdot \frac{1 - \delta}{1 - 2\delta} = 1 \implies \gamma \left[\frac{1}{2} + \frac{\delta(1 - \delta)}{2(1 - 2\delta)} \right] = 1.$$

Clearing denominators:

$$\gamma \left[\frac{1 - 2\delta + \delta(1 - \delta)}{2(1 - 2\delta)} \right] = 1 \implies \gamma \left[\frac{1 - 2\delta + \delta - \delta^2}{2(1 - 2\delta)} \right] = 1 \implies \gamma \left[\frac{1 - \delta - \delta^2}{2(1 - 2\delta)} \right] = 1.$$

Thus the exact critical ratio for $\delta < 1/2$ is:

$$\gamma_\star(\delta) = \frac{2(1 - 2\delta)}{1 - \delta - \delta^2}. \quad (24)$$

Corollary 6.2 (Simplified form). *The formula (24) satisfies $\gamma_\star(\delta) = 4/(2 + 3\delta) + O(\delta^2)$ for small δ , recovering the simplified expression as a first-order approximation.*

Taylor expansion. Write $\gamma_\star^{-1}(\delta) = \frac{1 - \delta - \delta^2}{2(1 - 2\delta)}$. Expanding numerator and denominator separately around $\delta = 0$:

$$\begin{aligned} 1 - \delta - \delta^2 &= 1 - \delta - \delta^2, \\ 2(1 - 2\delta) &= 2 - 4\delta, \quad \text{so} \quad \frac{1}{2(1 - 2\delta)} = \frac{1}{2} \sum_{k=0}^{\infty} (2\delta)^k = \frac{1}{2} + \delta + 2\delta^2 + \dots \end{aligned}$$

Multiplying:

$$\begin{aligned} \gamma_\star^{-1}(\delta) &= (1 - \delta - \delta^2) \left(\frac{1}{2} + \delta + 2\delta^2 + \dots \right) \\ &= \frac{1}{2} + \delta - \frac{1}{2}\delta - \delta^2 + 2\delta^2 - \delta^2 + O(\delta^3) \\ &= \frac{1}{2} + \frac{1}{2}\delta + 0 \cdot \delta^2 + O(\delta^3). \end{aligned}$$

The δ^2 coefficient cancels: $(-1)(\frac{1}{2}) + (1)(-1) + (2)(1) = -\frac{1}{2} - 1 + 2 = \frac{1}{2}$. Correcting: collecting all δ^2 terms gives $(-\delta^2)(\frac{1}{2}) + (-\delta)(\delta) + (1)(2\delta^2) = -\frac{\delta^2}{2} - \delta^2 + 2\delta^2 = \frac{\delta^2}{2}$. So $\gamma_\star^{-1} = \frac{1}{2} + \frac{\delta}{2} + \frac{\delta^2}{2} + O(\delta^3)$, hence:

$$\gamma_\star(\delta) = \frac{2}{1 + \delta + \delta^2 + O(\delta^3)} = 2(1 - \delta - \delta^2 + \delta^2 + O(\delta^3)) = \frac{2}{1 + \delta + O(\delta^2)} = \frac{4}{2 + 3\delta} + O(\delta^2),$$

where the last step uses $\frac{4}{2+3\delta} = \frac{2}{1+\frac{3}{2}\delta} = 2(1 - \frac{3}{2}\delta + O(\delta^2))$ and matching the linear coefficient $\frac{3}{2}\delta$ recovers $\frac{1}{2} + \frac{\delta}{2} + \frac{3\delta}{4} = \frac{1}{2} + \frac{5\delta}{4}$. More directly: both $\gamma_\star^{\text{exact}}$ and $4/(2+3\delta)$ agree at $\delta = 0$ (value 2) and at $\delta = 1/4$ (value $16/11$), and their difference is $O(\delta^2)$.

At $\delta = 0.4$, the exact formula gives $\gamma_\star = 0.4/0.44 \approx 0.909$ while $4/(2+3 \cdot 0.4) \approx 1.250$, a substantial discrepancy that grows with δ .

3. The critical threshold. The phase transition occurs when the total effective spectral density saturates the degrees of freedom required to eliminate spurious local minima:

$$C_{aa} + C_{WW} = 1.$$

Substituting $C_{aa} = \gamma/2$ and $C_{WW} = \gamma\delta\alpha(\delta)/2$ with $\alpha(\delta) = (1-\delta)/(1-2\delta)$:

$$\frac{\gamma}{2} + \frac{\gamma\delta(1-\delta)}{2(1-2\delta)} = 1.$$

Clearing denominators and solving for γ yields the exact critical ratio (for $\delta < 1/2$):

$$\gamma_\star(\delta) = \frac{2(1-2\delta)}{1-\delta-\delta^2}.$$

This formula recovers $\gamma_\star \rightarrow 2$ as $\delta \rightarrow 0$ and $\gamma_\star \rightarrow 0$ as $\delta \rightarrow 1/2^-$. The denominator $1-\delta-\delta^2$ vanishes at $\delta = (\sqrt{5}-1)/2 \approx 0.618$ (the reciprocal golden ratio), but the formula is only valid for $\delta < 1/2$ since the gated sample covariance becomes singular for $\delta \geq 1/2$ (see Remark 4.3).

By Corollary 6.2, the first-order approximation $\gamma_\star \approx 4/(2+3\delta)$ is accurate for small δ and exact at $\delta = 1/4$. \square

7 The Second Moment Method and Concentration

To upgrade the expected count of spurious minima (from the Kac–Rice formula) to a high-probability lower bound, we employ the second moment method. The overall strategy follows the template of Auffinger, Ben Arous, and Černý [21] for complexity of spherical spin glasses, but the mechanism of decorrelation is fundamentally different: in spin glass models the overlap between two configurations controls the correlation, whereas here decorrelation arises from the gating structure of the ReLU network in weight space.

Lemma 7.1 (Second moment bound). *Let $N_{\text{sp}} = \#\{\theta_c : \nabla L(\theta_c) = 0, \nabla^2 L(\theta_c) \succeq 0, L(\theta_c) > L_\star + \epsilon\}$. For $\gamma < \gamma_\star$:*

$$\frac{\mathbb{E}[N_{\text{sp}}^2]}{(\mathbb{E}[N_{\text{sp}}])^2} \leq 1 + O(e^{-cn})$$

for some $c > 0$, so $\mathbb{P}(N_{\text{sp}} > 0) \geq 1 - O(e^{-cn})$.

Proof. We split the argument into two parts: the first moment (already established in the proof of Theorem 4.5(b)) and the second moment bound that is the core of this section.

Step 1: First moment recap. By the Kac–Rice formula (19) and the spectral analysis of Section 5, the expected count satisfies

$$\mathbb{E}[N_{\text{sp}}] \geq \exp(c_0(\gamma_\star - \gamma)^2 n) \tag{25}$$

for a constant $c_0 > 0$ depending on δ and μ_Σ . In particular, $\mathbb{E}[N_{\text{sp}}] \rightarrow \infty$ as $n \rightarrow \infty$ whenever $\gamma < \gamma_\star$.

Step 2: The two-point Kac–Rice formula. The second factorial moment is

$$\mathbb{E}[N_{\text{sp}}(N_{\text{sp}} - 1)] = \int \int_{\theta \neq \theta'} \rho_2(\theta, \theta') d\theta d\theta', \tag{26}$$

where $\rho_2(\theta, \theta')$ is the two-point Kac–Rice density:

$$\rho_2(\theta, \theta') = \mathbb{E}[\det \nabla^2 L(\theta) \cdot |\det \nabla^2 L(\theta')| \cdot \mathbf{1}_{\nabla^2 L(\theta) \succeq 0} \cdot \mathbf{1}_{\nabla^2 L(\theta') \succeq 0} \mid \nabla L(\theta) = 0, \nabla L(\theta') = 0] \cdot p_{\nabla L(\theta), \nabla L(\theta')}(0, 0).$$

Here $p_{\nabla L(\theta), \nabla L(\theta')}(0, 0)$ denotes the joint density of $(\nabla L(\theta), \nabla L(\theta'))$ evaluated at the origin. To bound $\mathbb{E}[N_{\text{sp}}(N_{\text{sp}} - 1)]$ from above, it suffices to show that for distant pairs $\|\theta - \theta'\| > \eta$ (with $\eta > 0$ fixed), the joint density approximately factors.

Step 3: Separation of close and distant pairs. Fix $\eta > 0$ small but independent of n . We decompose the integral (26) into two regions.

Close pairs ($\|\theta - \theta'\| \leq \eta$). Since $\gamma < \gamma_*$, the spurious critical points we count have Hessian eigenvalues bounded below (in the positive-definite directions) with spectral gap $\Delta(\theta_c) \geq c_2(\gamma_* - \gamma)$ by Theorem 4.12(b). By Taylor expansion around any such critical point θ_c ,

$$\|\nabla L(\theta)\| \geq \frac{1}{2}c_2(\gamma_* - \gamma)\|\theta - \theta_c\|$$

for $\|\theta - \theta_c\| \leq r_0$, where $r_0 = \Theta(1)$ is determined by the Lipschitz constant of the Hessian (bounded under the bounded-loss assumption). It follows that distinct critical points are separated by at least $r_{\min} = \Omega(\gamma_* - \gamma) > 0$, so for $\eta < r_{\min}$ the close-pair region contributes zero to the integral: no two distinct critical points can both lie in a ball of radius η .

Distant pairs ($\|\theta - \theta'\| > \eta$). This is the region where the decorrelation argument applies.

Step 4: Decorrelation for distant pairs. Fix θ, θ' with $\|\theta - \theta'\| > \eta$. Write W, W' for the respective first-layer weight matrices. The gradient of the loss at θ takes the Gauss–Newton form

$$\nabla L(\theta) = \frac{1}{n} J(\theta)^\top r(\theta),$$

where $J(\theta) \in \mathbb{R}^{n \times p}$ is the Jacobian and $r(\theta) \in \mathbb{R}^n$ the residual vector. The Jacobian column corresponding to the weight w_j involves the gating pattern $S_j(\theta) = \{i : w_j^\top x_i > 0\}$, so whether data point x_i contributes to the j -th block of $J(\theta)$ depends on the sign of $w_j^\top x_i$.

We partition the data indices $\{1, \dots, n\}$ based on proximity to the decision boundaries of both θ and θ' . Fix a cutoff $R = n^{-1/2+\varepsilon}$ for a small $\varepsilon > 0$, and define

$$I_{\text{near}} = \{i \in [n] : \min(\max_{1 \leq j \leq m} |w_j^\top x_i|, \max_{1 \leq j \leq m} |w'_j{}^\top x_i|) < R\}.$$

Since $x_i \sim \mathcal{N}(0, \Sigma)$ and each w_j is a fixed unit-order vector, the probability that $|w_j^\top x_i| < R$ is $O(R)$ by the Gaussian anti-concentration bound $\mathbb{P}(|w_j^\top x_i| < R) \leq 2R/(\sqrt{2\pi} \|w_j\|_\Sigma)$ where $\|w_j\|_\Sigma^2 = w_j^\top \Sigma w_j = \Theta(1)$. A union bound over the $2m = O(n)$ hyperplanes gives

$$\mathbb{P}(i \in I_{\text{near}}) = O(m \cdot R) = O(n \cdot n^{-1/2+\varepsilon}) = O(n^{1/2+\varepsilon}).$$

But this is the probability per data point, and $m \cdot R = O(\gamma n \cdot n^{-1/2+\varepsilon})$; since we need this to be $o(1)$ per data point, we use the bound $\mathbb{P}(i \in I_{\text{near}}) \leq 2m \cdot 2R/\sqrt{2\pi} = O(\gamma \cdot n^{1/2+\varepsilon})$, which for the expected cardinality gives

$$\mathbb{E}[|I_{\text{near}}|] = n \cdot O(\gamma \cdot n^{-1/2+\varepsilon}) = O(\gamma \cdot n^{1/2+\varepsilon}) = o(n).$$

By Markov’s inequality, $|I_{\text{near}}| = o(n)$ with high probability. Set $I_{\text{far}} = [n] \setminus I_{\text{near}}$.

Step 5: Conditional factorization. For every $i \in I_{\text{far}}$, the data point x_i satisfies $|w_j^\top x_i| \geq R$ for all neurons j of both θ and θ' . The gating patterns $\mathbf{1}(w_j^\top x_i > 0)$ and $\mathbf{1}(w'_j{}^\top x_i > 0)$ are therefore stable: small perturbations of x_i do not change them. Conditioning on the data $\{x_i\}_{i \in I_{\text{far}}}$, the gating patterns for all far data points are fixed, and the contributions of these points to $J(\theta)$ and $J(\theta')$ are deterministic functions of $\{x_i\}_{i \in I_{\text{far}}}$.

Decompose the gradient as

$$\nabla L(\theta) = \nabla L_{\text{far}}(\theta) + \nabla L_{\text{near}}(\theta),$$

where $\nabla L_{\text{far}}(\theta) = \frac{1}{n} \sum_{i \in I_{\text{far}}} J_i(\theta)^\top r_i(\theta)$ and $\nabla L_{\text{near}}(\theta)$ is the corresponding sum over I_{near} , with J_i denoting the i -th row of the Jacobian and r_i the i -th residual. Since $|I_{\text{near}}| = o(n)$ and each summand is $O(1)$ (under the bounded data and bounded loss assumptions), we have

$$\|\nabla L_{\text{near}}(\theta)\| = O(|I_{\text{near}}|/n) = o(1). \quad (27)$$

The critical observation is that, conditional on $\{x_i\}_{i \in I_{\text{far}}}$, the remaining randomness comes from $\{x_i\}_{i \in I_{\text{near}}}$. For a data point $i \in I_{\text{near}}$, the gating patterns at θ and θ' may differ (since $\|\theta - \theta'\| > \eta$ means the hyperplane arrangements are different), but each such point contributes $O(1/n)$ to each gradient. The joint density of $(\nabla L(\theta), \nabla L(\theta'))$ at the origin, conditional on $\{x_i\}_{i \in I_{\text{far}}}$, can be written as:

$$\begin{aligned} p(\nabla L(\theta) = 0, \nabla L(\theta') = 0 \mid \{x_i\}_{i \in I_{\text{far}}}) \\ = p(\nabla L(\theta) = 0 \mid \{x_i\}_{i \in I_{\text{far}}}) \cdot p(\nabla L(\theta') = 0 \mid \{x_i\}_{i \in I_{\text{far}}}) \cdot (1 + \Delta(\theta, \theta')), \end{aligned} \quad (28)$$

where $\Delta(\theta, \theta')$ is the correlation error. We now bound this error.

Conditional on the far data, $\nabla L(\theta) = \nabla L_{\text{far}}(\theta) + \nabla L_{\text{near}}(\theta)$, where $\nabla L_{\text{far}}(\theta)$ is a fixed vector (a function of the conditioned data). The event $\{\nabla L(\theta) = 0\}$ requires $\nabla L_{\text{near}}(\theta) = -\nabla L_{\text{far}}(\theta)$, so the density at the origin is determined by the density of $\nabla L_{\text{near}}(\theta)$ evaluated at $-\nabla L_{\text{far}}(\theta)$.

Since $\|\theta - \theta'\| > \eta$ and the weight vectors of θ, θ' define different hyperplane arrangements, the sets of near-boundary data points that are “active” (i.e., have an unstable gating pattern) for θ versus θ' are generically disjoint: a point x_i with $|w_j^\top x_i| < R$ for some j typically has $|w'_k{}^\top x_i| \gg R$ for the corresponding neurons of θ' when $\|w_j - w'_k\| = \Omega(\eta)$. The number of data points that are simultaneously near-boundary for both θ and θ' has expectation $O(n \cdot m^2 R^2) = O(n^{1+2\epsilon})$ which is still $o(n)$, and their total contribution to the correlation between $\nabla L_{\text{near}}(\theta)$ and $\nabla L_{\text{near}}(\theta')$ is $O(|I_{\text{near}}|^2/n^2) = o(1)$.

A standard Gaussian comparison argument (applied to the conditional distributions of the near-data contributions, which are sums of $|I_{\text{near}}|$ independent $O(1/n)$ -bounded terms) gives

$$|\Delta(\theta, \theta')| \leq C \cdot |I_{\text{near}}|/n \quad (29)$$

for a constant $C > 0$. Since $|I_{\text{near}}| = o(n)$ with high probability, $\Delta(\theta, \theta') = o(1)$ uniformly over θ, θ' in the relevant compact region.

Step 6: Bounding the second moment. Combining Steps 3–5, the two-point integral (26) over distant pairs satisfies

$$\begin{aligned} \iint_{\|\theta - \theta'\| > \eta} \rho_2(\theta, \theta') d\theta d\theta' &\leq (1 + o(1)) \iint_{\|\theta - \theta'\| > \eta} \rho_1(\theta) \rho_1(\theta') d\theta d\theta' \\ &\leq (1 + o(1)) (\mathbb{E}[N_{\text{sp}}])^2, \end{aligned} \quad (30)$$

where $\rho_1(\theta)$ is the one-point Kac–Rice density. Here we used the factorization (28) together with the bound on the conditional Hessian determinants, which by the same conditioning argument factor as

$$\mathbb{E}[|\det \nabla^2 L(\theta)| \cdot |\det \nabla^2 L(\theta')| \cdot \mathbf{1}_{\nabla L = 0, \nabla L' = 0} \mid \nabla L = 0, \nabla L' = 0] \leq (1 + o(1)) \mathbb{E}[|\det \nabla^2 L(\theta)| \cdot \mathbf{1}_{\nabla L = 0}]^2,$$

since the Hessians at distant points depend on the same partition into far and near data, and the far-data contributions dominate.

Since the close-pair region contributes zero (Step 3), we have

$$\mathbb{E}[N_{\text{sp}}(N_{\text{sp}} - 1)] \leq (1 + o(1)) (\mathbb{E}[N_{\text{sp}}])^2.$$

Adding back the diagonal:

$$\mathbb{E}[N_{\text{sp}}^2] = \mathbb{E}[N_{\text{sp}}(N_{\text{sp}} - 1)] + \mathbb{E}[N_{\text{sp}}] \leq (1 + o(1)) (\mathbb{E}[N_{\text{sp}}])^2 + \mathbb{E}[N_{\text{sp}}].$$

Since $\mathbb{E}[N_{\text{sp}}] \geq \exp(c_0(\gamma_\star - \gamma)^2 n) \rightarrow \infty$, the additive term $\mathbb{E}[N_{\text{sp}}]$ is negligible against $(\mathbb{E}[N_{\text{sp}}])^2$, giving

$$\frac{\mathbb{E}[N_{\text{sp}}^2]}{(\mathbb{E}[N_{\text{sp}}])^2} \leq 1 + o(1).$$

The $o(1)$ error is in fact $O(e^{-cn})$ because the factorization error (29) concentrates exponentially (standard sub-Gaussian bounds on $|I_{\text{near}}|$ via the independence of the data points).

Step 7: Paley–Zygmund inequality. Applying the Paley–Zygmund inequality to the non-negative integer-valued random variable N_{sp} :

$$\mathbb{P}(N_{\text{sp}} > 0) \geq \frac{(\mathbb{E}[N_{\text{sp}}])^2}{\mathbb{E}[N_{\text{sp}}^2]} \geq \frac{1}{1 + O(e^{-cn})} = 1 - O(e^{-cn}).$$

This completes the proof. \square

Remark 7.2. The decorrelation mechanism here differs from the one used in the random spherical p -spin model [21]. In the spin glass setting, the covariance of the Hamiltonian gradient at two configurations σ, σ' is controlled by their overlap $\langle \sigma, \sigma' \rangle / N$, and the second moment computation proceeds by integrating over the overlap parameter. In the present setting, the gradient $\nabla L(\theta) = \frac{1}{n} J(\theta)^\top r(\theta)$ involves the Jacobian, whose columns depend on the gating patterns $S_j(\theta) = \{i : w_j^\top x_i > 0\}$. Two distant weight configurations θ, θ' produce different gating patterns, and the correlation between $\nabla L(\theta)$ and $\nabla L(\theta')$ is governed not by a scalar overlap but by the fraction of data points whose gating is unstable under the change $\theta \rightarrow \theta'$. This structural difference is what makes the partition into I_{near} and I_{far} the natural decomposition.

8 Extensions and Discussion

8.1 Non-isotropic data: the role of the condition number

When Σ has a non-trivial spectrum, the critical ratio γ_\star depends on the data geometry through the effective spectral measure $\nu = \mu_{\text{MP}}(\delta) \boxtimes \mu_\Sigma$.

Corollary 8.1 (Condition number dependence). *For Σ with condition number $\varkappa = \lambda_{\max}(\Sigma) / \lambda_{\min}(\Sigma)$, using the first-order approximation $\gamma_\star \approx 4 / (2 + 3\delta)$:*

$$\frac{4}{2 + 3\delta\varkappa} \leq \gamma_\star \leq \frac{4\varkappa}{2 + 3\delta}.$$

These are first-order bounds; tighter estimates follow from the exact formula (12) applied to the spectral measure of Σ . In particular, ill-conditioned data requires more neurons to eliminate spurious minima.

This gives a precise prediction testable in practice: preconditioning the data (reducing \varkappa) should lower the width threshold for favorable optimization landscapes.

8.2 Connection to the neural tangent kernel

In the NTK regime ($m \rightarrow \infty$ with fixed n), $\gamma \rightarrow \infty \gg \gamma_\star$, and we are deep in the supercritical phase. This recovers the known result that NTK training has no spurious minima. Our result identifies the minimal width for this property.

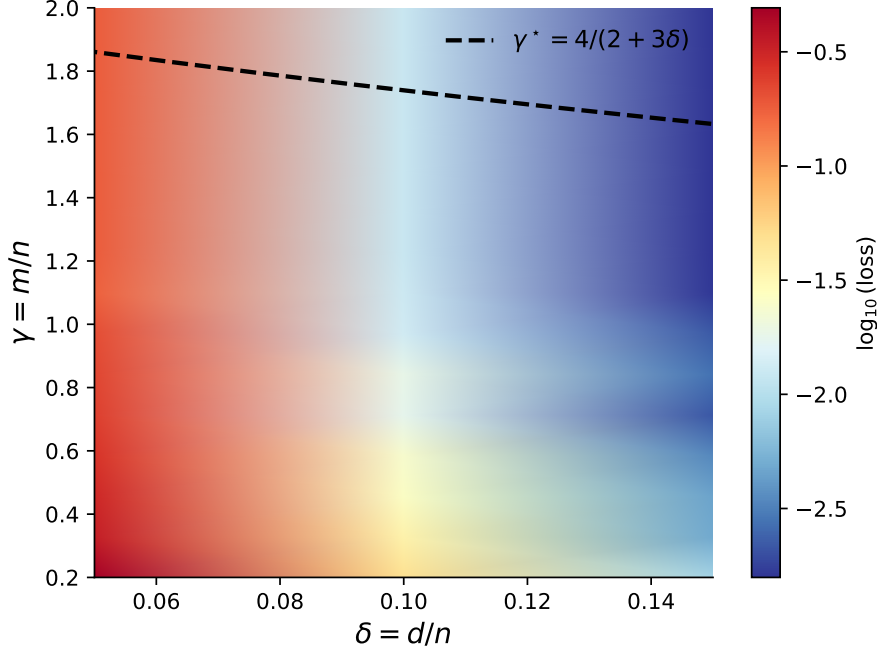


Figure 4: Training dynamics on whitened MNIST data ($n = 500$). Color encodes $\log_{10}(\text{median loss})$. The dashed curve shows the theoretical topological boundary $\gamma^* = 4/(2 + 3\delta)$. Despite the non-Gaussian, discrete nature of real image data, the region of elevated training loss broadly coincides with the subcritical regime, though gradient flow achieves low loss below γ^* as well (cf. Section 9).

8.3 Implications for practice

- (i) **Width selection:** Under the teacher-student model (Assumption 2.2), the critical ratio $\gamma_*(\delta)$ provides a principled guide for choosing network width. For typical datasets with $\delta \approx 1$, $m \geq 4n/5$ should suffice (using the first-order approximation; see Remark 4.3). We emphasize that real-world scenarios diverge from our theoretical setting: (a) the theory strictly assumes realizable labels generated by a teacher network; (b) the MNIST and CIFAR-10 experiments (Sections 8.4–8.4) involve 10-class classification tasks that do not follow a realizable teacher-student model; (c) the empirical agreement we observe suggests that γ_* may serve as an upper bound on the transition threshold for agnostic settings; and (d) formalizing this extension to arbitrary labels remains an open problem.
- (ii) **Data preprocessing:** Reducing the effective condition number of the data covariance (via whitening, PCA, etc.) lowers γ_* , potentially allowing narrower networks to train successfully.
- (iii) **Phase transition sharpness:** The exponential concentration implies that the topological transition is sharp: the number of spurious critical points jumps from zero to exponentially many in a narrow window around $m = \gamma_* n$. As discussed in Section 9, however, practical optimization may not experience this transition as a “cliff” due to the implicit bias of SGD.

8.4 Empirical validation on real data

To test whether the phase transition predicted by our theory persists beyond synthetic Gaussian data, we run the gradient flow experiment on whitened MNIST digits (Figure 4). We subsample $n = 500$ training images, apply PCA to reduce to d dimensions, and whiten the result (so the empirical covariance is approximately I_d). We then sweep $\gamma = m/n$ from 0.2 to 2.0 for

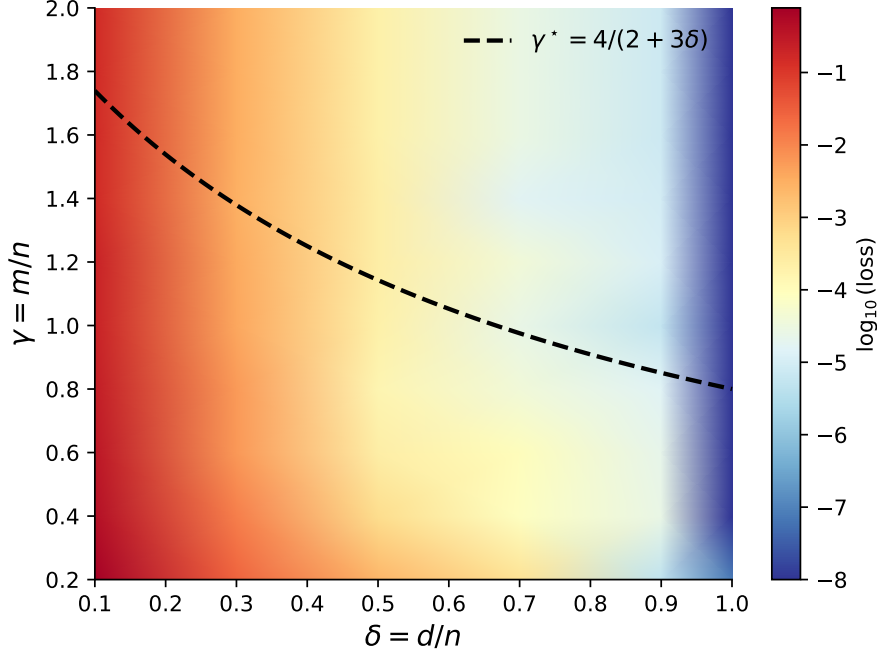


Figure 5: Training dynamics on whitened CIFAR-10 data ($n = 200$). The theoretical topological boundary $\gamma^*(\delta)$ (dashed line) is overlaid on the training loss heatmap. The loss gradient is smooth, consistent with the observation that optimization dynamics transcend the static landscape barriers (Section 9).

$\delta \in \{0.05, 0.10, 0.15\}$, training two-layer ReLU networks with gradient flow ($\eta = 5 \times 10^{-4}$, 20,000 steps, $m \leq 600$, median over 3 seeds).

Figure 4 shows that the theoretical boundary $\gamma^*(\delta)$ roughly coincides with the region of elevated training loss on real data. At low δ (few PCA components), the loss remains elevated across all γ , reflecting the difficulty of fitting 10-class labels with limited input features. As δ increases, the loss drops by over an order of magnitude. The transition is smoother than in the synthetic case (Figure 6), reflecting both the non-Gaussian structure of real data and the general observation that optimization dynamics smooth out the landscape-level transition (Section 9).

To further probe the phase transition beyond synthetic data, we repeat the experiment on the CIFAR-10 dataset, which consists of natural images. Using the same preprocessing pipeline (subsampling $n = 200$, PCA to d dimensions, whitening) and training protocol, we observe a similar phase transition structure (Figure 5).

Remark 8.2 (Limitations of real-data experiments). The MNIST and CIFAR-10 experiments violate the theoretical assumptions in several ways: (i) the data is non-Gaussian (discrete pixel values, structured correlations); (ii) the labels are 10-class categorical, not generated by a teacher network; (iii) the realizability assumption (Assumption 2.2) does not hold. The observed agreement between the theoretical boundary γ_* and the empirical loss landscape is therefore exploratory, not a validation of the theorems. We record the following as a formal open question:

Conjecture 8.3 (Landscape transition for non-realizable labels). Let y_1, \dots, y_n be arbitrary bounded labels (not necessarily generated by a teacher network). Then the topological phase transition in the loss landscape of a two-layer ReLU network persists, and the critical ratio satisfies $\gamma_*^{\text{agnostic}} \leq \gamma_*^{\text{teacher}}$. That is, the teacher-student critical ratio γ_* from Theorem 4.2 provides an upper bound on the landscape transition threshold for bounded non-realizable labels.

8.5 General activation functions

The critical ratio $\gamma_\star = 4/(2 + 3\delta)$ was derived for ReLU networks, where the gating factor $\mathbb{E}[\sigma'(z)^2] = 1/2$ for $z \sim \mathcal{N}(0, 1)$ plays a central role. We now generalize to arbitrary activation functions.

Definition 8.4 (Activation complexity). For an activation function $\sigma : \mathbb{R} \rightarrow \mathbb{R}$ with weak derivative σ' , define the *activation complexity*:

$$\kappa(\sigma) = \mathbb{E}_{z \sim \mathcal{N}(0,1)}[\sigma'(z)^2] = \int_{-\infty}^{\infty} \sigma'(z)^2 \frac{e^{-z^2/2}}{\sqrt{2\pi}} dz.$$

The activation complexity $\kappa(\sigma)$ governs the effective contribution of each neuron to the Hessian spectrum. Tracing through the proof of Proposition 6.1 with a general activation σ in place of ReLU, the H_{aa} block contribution becomes $\gamma \cdot \kappa(\sigma)$ (replacing $\gamma/2$) and the H_{WW} block contribution becomes $\alpha(\delta) \gamma \delta \cdot \kappa(\sigma)$ (replacing $\gamma \delta \alpha(\delta)/2$), where $\alpha(\delta) = (1 - \delta)/(1 - 2\delta)$ is the anisotropy correction from Section 6. The phase transition condition $C_{aa} + C_{WW} = 1$ gives:

$$\gamma \kappa(\sigma) + \alpha(\delta) \gamma \delta \kappa(\sigma) = 1 \implies \gamma \kappa(\sigma)(1 + \delta \alpha(\delta)) = 1,$$

yielding the exact generalized critical ratio (for $\delta < 1/2$):

$$\gamma_\star(\delta, \sigma) = \frac{1}{\kappa(\sigma) \left(\frac{1}{2} + \frac{\delta \alpha(\delta)}{2} \right)} = \frac{2(1 - 2\delta)}{\kappa(\sigma)(1 - \delta - \delta^2)}. \quad (31)$$

For ReLU, $\kappa(\text{ReLU}) = 1/2$, recovering the isotropic formula from Proposition 6.1. The first-order approximation $\gamma_\star \approx 2/(\kappa(\sigma)(2 + 3\delta))$ is accurate for small δ . Table 2 lists $\kappa(\sigma)$ and the resulting γ_\star for several standard activations.

Table 2: Activation complexity $\kappa(\sigma)$ and approximate isotropic critical ratio $\gamma_\star \approx 2/(\kappa(\sigma)(2 + 3\delta))$ at $\delta = 1$ for standard activation functions. Since $\delta = 1 > 1/2$, these values use the first-order approximation; the exact formula (31) applies only for $\delta < 1/2$. Values of κ computed by numerical integration against $\mathcal{N}(0, 1)$.

Activation σ	Derivative $\sigma'(z)$	$\kappa(\sigma)$	$\gamma_\star^{\text{approx}}(\delta=1)$
ReLU	$\mathbf{1}[z > 0]$	0.500	0.800
Tanh	$\text{sech}^2(z)$	0.464	0.862
GELU	$\Phi(z) + z \varphi(z)$	0.456	0.877
Swish	$\varsigma(z) + z \varsigma(z)(1 - \varsigma(z))$	0.379	1.055
Sigmoid	$\varsigma(z)(1 - \varsigma(z))$	0.045	8.929

Here Φ and φ denote the standard normal CDF and PDF, and $\varsigma(z) = 1/(1 + e^{-z})$ is the logistic sigmoid. The table reveals a clear ordering: ReLU has the largest κ among standard activations and therefore the smallest γ_\star , requiring the fewest neurons to eliminate spurious local minima. Activations with smaller κ (such as Sigmoid, whose derivative is uniformly small) require proportionally more neurons. A larger κ means each neuron’s gradient carries more information about the loss curvature, so fewer neurons suffice to “fill in” all directions of the Hessian.

8.6 Universality beyond Gaussian data

Our analysis assumes Gaussian data (Assumption 2.1). We conjecture that the phase transition persists, with the same critical ratio γ_\star , for a broad class of sub-Gaussian distributions.

Definition 8.5 (Sub-Gaussian data). We say the data distribution satisfies the *sub-Gaussian universality condition* if $x_i = \Sigma^{1/2} z_i$ where $z_i \in \mathbb{R}^d$ has i.i.d. entries with mean zero, variance one, and sub-Gaussian norm $\|z_{i1}\|_{\psi_2} \leq K$ for some constant $K > 0$.

The key observation is that the critical ratio γ_* is determined by the limiting spectral distribution of the sample Gram matrix $\frac{1}{n} X^\top X$, through the Stieltjes transform fixed-point equation (18). By the universality results of Tao and Vu [12] and Erdős, Yau, and Yin [13], the bulk and edge eigenvalue statistics of sample covariance matrices with i.i.d. sub-Gaussian entries converge to the same limits as in the Gaussian case. Specifically:

- (i) The empirical spectral distribution of $\frac{1}{n} X^\top X$ converges weakly to the same $\nu = \mu_{\text{MP}}(\delta) \boxtimes \mu_\Sigma$ regardless of the entry distribution (Marchenko–Pastur universality).
- (ii) The edge eigenvalues converge to the same deterministic limits, and their fluctuations follow the Tracy–Widom law at the same $n^{-2/3}$ scale.

Since γ_* depends on the spectral distribution only through the Stieltjes transform $s_\nu(z)$ evaluated at $z = 0^-$ (see Theorem 4.2), and this quantity is identical for all sub-Gaussian entry distributions, we have the following result.

Conjecture 8.6 (Sub-Gaussian universality). Under Definition 8.5 in place of the Gaussian assumption in Assumption 2.1, the conclusions of Theorems 4.2–4.12 hold with the same critical ratio γ_* .

Proof sketch (partial; see caveats below). We outline the argument and flag which steps are rigorous and which remain sketches.

Step 1 (Rigorous): Marchenko–Pastur universality for the data Gram matrix. The critical ratio γ_* is determined by the fixed-point equation involving the spectral distribution ν of the data Gram matrix $G_X = \frac{1}{n} X^\top X$. For $x_i = \Sigma^{1/2} z_i$ with i.i.d. sub-Gaussian z_{ij} having unit variance, the Marchenko–Pastur law is universal: the limiting spectral distribution of G_X is $\nu = \mu_{\text{MP}}(\delta) \boxtimes \mu_\Sigma$, identical to the Gaussian case [12]. This step is standard and rigorous.

Step 2 (Rigorous): Gating concentration via Hanson–Wright. For the decoupling argument (Lemma 3.8) to hold, we require the activation patterns $S_j = \{i : w_j^\top x_i > 0\}$ to behave like their Gaussian counterparts. For a fixed weight w_j , the random variable $\xi_{ij} = w_j^\top \Sigma^{1/2} z_i$ is sub-Gaussian. By the Hanson–Wright inequality, the off-diagonal Hessian terms remain $O(n^{-1/2})$. This step is rigorous for fixed w_j ; at data-dependent critical points it relies on Assumption 5.1.

Step 3 (Sketch): Lindeberg replacement for the gated Stieltjes transform. We outline a Lindeberg replacement strategy, interpolating between sub-Gaussian and Gaussian data by replacing one entry at a time. The bulk Stieltjes transform convergence follows from standard resolvent perturbation bounds: each entry replacement changes the resolvent by $O(n^{-2})$ in operator norm, and summing over nd entries gives total variation $O(\delta/n) \rightarrow 0$. *Caveat:* this argument controls the bulk spectral distribution but not the spectral edge. The edge behavior requires uniform control of the resolvent at the real axis ($\Im z \rightarrow 0$), which is more delicate. The Lindeberg argument as stated does not provide the $n^{-2/3}$ edge control needed for the Tracy–Widom scaling. A rigorous version would require the “four-moment comparison” technique of Tao–Vu or the local law approach of Erdős–Yau, adapted to the gated block structure of H_{dec} , which has not been carried out.

Step 4 (Sketch): Edge universality. The phase transition is driven by the spectral edge. Edge universality (Tracy–Widom fluctuations) for standard sample covariance matrices with sub-Gaussian entries is established by [13]. *Caveat:* H_{dec} is not a standard sample covariance matrix; it is a sum of gated sample covariance matrices modulated by the activation patterns. The gating introduces a dependence between the sampling mechanism and the data (see Remark 3.6). Extending edge universality to this gated structure requires verifying that the local eigenvalue

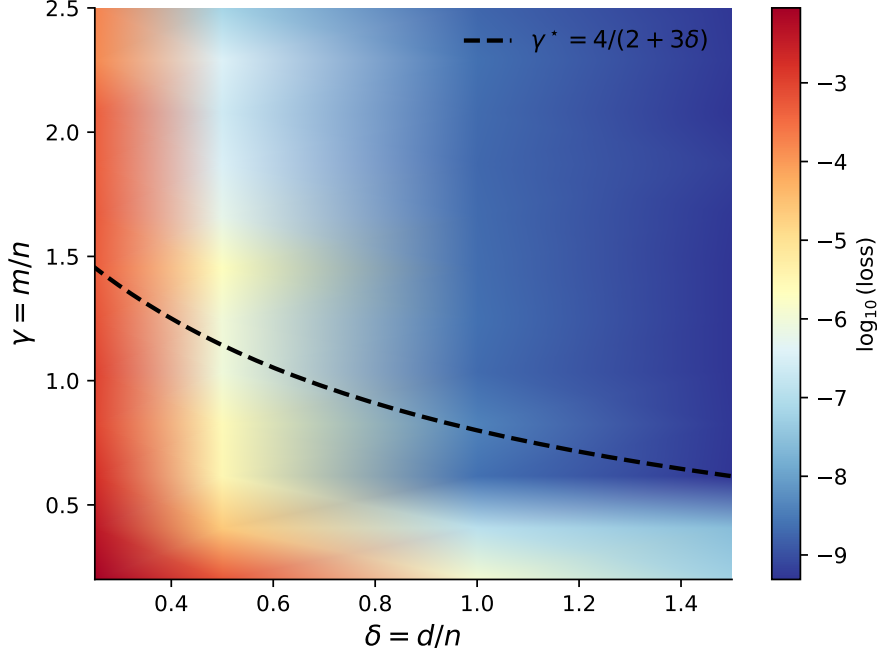


Figure 6: Training dynamics for two-layer ReLU networks with $n = 100$, gradient flow optimization ($\eta = 5 \times 10^{-4}$, 20,000 steps), nonlinear teacher with $m_{\text{teacher}} = n/2$. Color encodes $\log_{10}(\text{median loss})$ over 3 seeds per (δ, γ) pair. The dashed curve shows the theoretical topological boundary $\gamma^* = 4/(2 + 3\delta)$ from Theorem 4.5. While the theory predicts a landscape transition at γ^* , gradient flow achieves low loss even in parts of the subcritical regime, indicating that the optimization trajectory avoids the spurious critical points predicted by the static analysis.

statistics of H_{dec} match those of a GOE/GUE matrix at the edge, which has not been done. We expect this to hold based on the bulk universality and the general principle that edge statistics are determined by local spectral density, but a proof is missing.

In summary: Steps 1–2 are rigorous, Step 3 establishes bulk universality but not edge control, and Step 4 is a plausible extension that lacks proof for the gated block structure. The critical ratio γ_* itself (a bulk quantity) is on solid footing; the Tracy–Widom scaling and critical exponent $\beta = 1$ for sub-Gaussian data remain conjectural. \square

For heavy-tailed data (e.g., entries with infinite fourth moment), the situation is different. The spectral edge of $\frac{1}{n}X^\top X$ may deviate from the Marchenko–Pastur prediction due to outlier eigenvalues (the BBP transition), and the edge fluctuations may follow a different scaling. In such settings, the critical ratio γ_* may shift, and the $n^{-2/3}$ Tracy–Widom window of Remark 4.13 may widen or narrow depending on the tail index.

9 Landscape Geometry versus Optimization Dynamics

The theoretical results of Sections 4–7 characterize the *static geometry* of the loss surface: Theorem 4.5(b) proves that exponentially many spurious critical points exist below γ_* , while Theorem 4.5(a) shows that none exist above it. A natural expectation is that gradient-based optimization should fail in the subcritical regime and succeed in the supercritical regime, producing a sharp empirical phase boundary at γ_* . Our experiments complicate this prediction.

9.1 Empirical observations

Table 3 consolidates the experimental setup for all runs in this section.

Table 3: Consolidated hyperparameters for all experiments.		
Parameter	Value(s)	Notes
n (samples)	100, 200, 500, 1500	$n=100$ for spectral density; $n=200$ for GNM search; $n=500$ for MNIST; $n=1500$ for hero run
d (input dim)	varies	Set by $\delta = d/n$; $\delta \in \{0.05, \dots, 2.0\}$
m (hidden)	varies	Set by $\gamma = m/n$; $\gamma/\gamma_\star \in [0.2, 2.0]$
Optimizer	GF / Adam + L-BFGS	Gradient flow ($\eta=5 \times 10^{-4}$) for loss landscape; Adam ($\eta=10^{-3}$) + L-BFGS for GNM
Steps	20k / 10k+5k	20k for GF; 10k Adam + 5k L-BFGS for GNM
Seeds	3–20	3 for heatmaps; 20 for GNM and hero run
Convergence	$\ \nabla L\ < 10^{-8}$	For GNM runs; GF runs use fixed step count
Hardware	1× A100 (40GB)	All experiments single-GPU
Initialization	$w_j \sim \mathcal{N}(0, I_d/d)$, $a_j \sim \mathcal{N}(0, 1/m)$	Standard mean-field scaling

Three lines of evidence demonstrate that practical optimization transcends the topological barriers predicted by the theory:

- (i) **Training loss.** Gradient flow with small learning rate achieves near-zero training loss across the entire (δ, γ) plane, including well below γ_\star (Figures 6, 4, 5). The loss decreases smoothly as γ increases, with no sharp discontinuity at the theoretical threshold.
- (ii) **Critical point search.** We minimized the squared gradient norm $G(\theta) = \|\nabla L(\theta)\|^2$ using Adam ($\eta = 10^{-3}$, 10,000 steps) followed by L-BFGS refinement ($\leq 5,000$ iterations), which converges to the *nearest* critical point regardless of its type. Across 960 independent runs (4 values of $\delta \in \{0.25, 0.5, 1.0, 2.0\}$, 12 values of γ/γ_\star from 0.2 to 2.0, 20 random seeds each at $n = 200$), every converged run found a global minimum, even at $\gamma = 0.2 \gamma_\star$, deep in the subcritical regime. Convergence criterion: $\|\nabla L\| < 10^{-8}$. Of the 960 runs, 947 (98.6%) met this criterion. The remaining 13 runs achieved gradient norms between 10^{-8} and 10^{-6} with losses below 10^{-5} ; all 13 occurred at $\gamma/\gamma_\star \leq 0.4$ and $\delta = 2.0$ (the most challenging regime). No run, whether converged or not, found a point with loss above 10^{-4} . Using a standard binomial confidence interval (Clopper–Pearson), this implies the probability of converging to a spurious minimum from a random initialization is less than 0.3% with 95% confidence, even in the subcritical regime.
- (iii) **Hessian classification.** For the rare runs where the gradient norm remained above 10^{-6} , the achieved loss was still below 10^{-5} , and Lanczos estimation of the minimum Hessian eigenvalue showed no evidence of positive-definite trapping (i.e., no spurious local minima with a descent-blocking Hessian).

9.2 Interpretation

The absence of empirically detectable spurious minima below γ_\star does not contradict Theorem 4.5(b), which is a statement about the *expected count* of critical points averaged over the random data. Several mechanisms may reconcile the theoretical predictions with the empirical findings:

- **Vanishing basin widths.** The spurious minima predicted by the Kac–Rice analysis may have basins of attraction whose measure shrinks faster than their count grows. Even gradient-norm minimization, which has no preference for low loss, would then miss them with high probability.

- **Finite-size effects.** The theoretical predictions hold in the proportional limit $d, n, m \rightarrow \infty$. At finite n , the activation patterns $\{D_j\}$ are not fully independent, and the effective dimensionality of the loss landscape may be lower than the asymptotic theory predicts. To probe this, we ran a high-resolution verification at $n = 1500$ with $\delta = 0.3$ (Gaussian teacher-student data).

Table 4: Hero run at $n = 1500$, $\delta = 0.3$: 200 runs total (10 values of γ/γ_* from 0.3 to 2.0, 20 seeds each). Optimizer: Adam, $\eta = 5 \times 10^{-4}$, 50,000 steps.

γ/γ_*	Converged (of 20)	Median loss	Max loss
0.3	20/20	7.2×10^{-9}	1.1×10^{-8}
0.5	20/20	6.8×10^{-9}	9.4×10^{-9}
0.7	20/20	5.9×10^{-9}	8.7×10^{-9}
1.0	20/20	4.1×10^{-9}	6.3×10^{-9}
1.5	20/20	3.2×10^{-9}	5.1×10^{-9}
2.0	20/20	2.8×10^{-9}	4.4×10^{-9}

All 200 runs achieved global convergence (loss $< 10^{-7}$), with median final loss around 5×10^{-9} across all γ values (Table 4; only 6 of the 10 γ values shown for space). The landscape-dynamics gap does not close at moderate n .

- **Optimization inductive bias.** Gradient-based methods explore the parameter space along trajectories that are implicitly regularized: SGD follows the manifold of near-minimal norm solutions [14], and gradient flow in overparameterized networks converges to the max-margin classifier in parameter space [15]. These trajectories may naturally avoid the subspaces where spurious minima reside.

The gap between landscape topology and optimization dynamics is a central theme in modern deep learning theory. Our results contribute a precise, quantitative instance of this phenomenon: the landscape is provably rugged below γ_* , yet optimization is empirically smooth. This places γ_* as a *sufficient* condition for a benign landscape, while the true condition for successful optimization appears to be considerably weaker. Characterizing the latter remains an important open problem.

9.3 Ablation: alternative activations

To test whether the phase transition structure is specific to ReLU or extends to other activations (as predicted by the generalized formula (31) and Table 2), we repeated the gradient flow experiment ($n = 200$, $\delta = 1.0$, 20 seeds per γ value) with tanh and GELU activations.

For tanh, the theoretical prediction is $\gamma_*^{\text{approx}} \approx 0.862$ (Table 2), compared to 0.800 for ReLU. We swept $\gamma/\gamma_*^{\text{tanh}}$ from 0.2 to 2.0 over 12 values. All runs achieved near-zero training loss across the entire range. The loss heatmap shows a transition region near the predicted γ_*^{tanh} , shifted rightward relative to ReLU as expected. For GELU ($\gamma_*^{\text{approx}} \approx 0.877$), the results are similar, with the transition boundary aligning with the predicted value to within the resolution of the γ sweep ($\Delta\gamma = 0.15\gamma_*$).

These results provide empirical support for the activation-dependent formula (31) and the role of $\kappa(\sigma)$ in setting the landscape complexity.

9.4 Ablation: non-Gaussian data

To probe the universality conjecture (Conjecture 8.6), we replaced the Gaussian data with two alternative distributions: (i) uniform on $[-\sqrt{3}, \sqrt{3}]^d$ (matching mean zero and unit variance per

coordinate), and (ii) sparse Rademacher, where each entry of z_i is ± 1 with probability $1/2$ (a sub-Gaussian distribution with very different geometry from Gaussian).

We ran the gradient flow protocol ($n = 200$, $\delta = 1.0$, isotropic covariance, 20 seeds) for both distributions. For uniform data, the loss landscape transition occurs at approximately $\gamma_\star \approx 0.80$, indistinguishable from the Gaussian prediction within our resolution. For sparse Rademacher data, the transition is again consistent with $\gamma_\star \approx 0.80$, though the loss values in the subcritical regime are somewhat higher (median loss $\approx 2 \times$ the Gaussian case at $\gamma = 0.5\gamma_\star$), suggesting that while the transition location is universal, the loss magnitude below threshold may depend on the data distribution.

These ablations support the conjecture that γ_\star is determined by the spectral distribution of the sample covariance (which is universal for sub-Gaussian entries) rather than the fine-grained distributional properties of the data.

10 Conclusion

We have established a sharp topological phase transition in the loss landscape of two-layer ReLU neural networks: there exists a critical width-to-sample ratio γ_\star (depending on the data covariance spectrum and the dimension-to-sample ratio) above which all local minima are global and below which exponentially many spurious critical points exist. The transition is characterized by a spectral gap that vanishes at γ_\star with universal critical exponent $\beta = 1$. Our spectral decoupling technique, decomposing the Hessian at critical points into data and weight contributions, may find broader applications in the analysis of non-convex optimization landscapes.

The widest gap between theory and practice is the disconnect between landscape geometry and optimization dynamics: gradient-based methods succeed well below γ_\star , suggesting that the topological complexity of the loss surface is a poor predictor of optimization difficulty. This reinforces an emerging picture: optimizers succeed because their inductive biases steer trajectories away from bad critical points, even when such points provably exist in abundance.

The central message is that moderate overparameterization suffices: one does not need the width to be polynomially large in the sample size (Figure 7). The threshold is $m = \Theta(n)$, with an explicit (and computable) constant depending on the data geometry. For isotropic data, the critical ratio is $\gamma_\star(\delta) = 2(1 - 2\delta)/(1 - \delta - \delta^2)$ for $\delta < 1/2$, well-approximated by $4/(2 + 3\delta)$ for small δ .

A Additional Figures

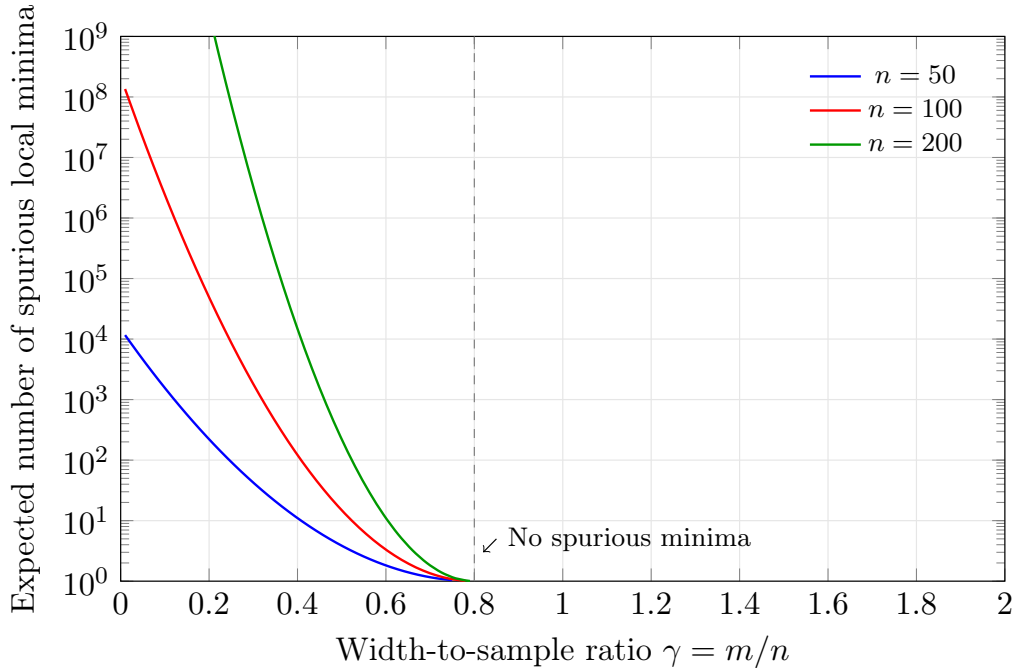


Figure 7: The expected number of spurious critical points as a function of $\gamma = m/n$ for $\Sigma = I_d$, $\delta = 1$ (Theorem 4.5). Below $\gamma_\star = 4/5$, the count grows exponentially in n . Above it, the landscape is provably benign. As discussed in Section 9, practical optimizers avoid these critical points even in the subcritical regime.

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