Local Coverage Governs Memorization in Diffusion Models

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Abstract

Memorization in diffusion models is often treated as a global property of the model or dataset. In practice, however, a single diffusion model can simultaneously generate both memorized and novel samples. Which training samples are most likely to be memorized? In this work, we show that memorization is governed by *local data coverage*. Leveraging the connection between diffusion models and kernel density estimation (KDE), we derive a theoretical criterion that predicts whether a point is memorized based on the density of training data in its neighborhood and the size of the training dataset. In the high-dimensional limit, this leads to a sharp, local transition: regions of low coverage are dominated by isolated training samples, which are memorized, while dense regions support interpolation and generalization. We validate these predictions empirically, showing that memorization increases with local sparsity and that diffusion models exhibit a coexistence of memorized and novel samples within the same model. Extending this framework to multiclass settings, we further show that classes with higher intra-class diversity (and thus lower local coverage) are more strongly memorized. Our results provide a local view of memorization in diffusion models, explaining when and where memorization occurs in terms of data geometry.

1. Introduction

How does memorization arise *locally* in diffusion models? Diffusion models are powerful generative methods capable of learning complex high-dimensional data distributions from finite datasets. Yet when trained with limited data, they may reproduce training examples rather than generate novel samples, a phenomenon commonly referred to as *memorization* (Somepalli et al., 2022; Carlini et al., 2023;

Kadkhodaie et al., 2024). Understanding the mechanisms that drive memorization is important because memorized samples can lead to unintended data leakage, privacy violations, and the release of copyrighted or sensitive training data.

Theoretical work has begun to analyze memorization in linear models (Maillard & Goldt, 2026; Wang et al., 2026), in random feature models (George et al., 2026; Bonnaire et al., 2026), and by exploiting the close connection between diffusion models and kernel density estimation (KDE) (Pidstrigach, 2022; Pham et al., 2024; Ambrogioni, 2024; Biroli & Mézard, 2024; Achilli et al., 2025a; Li et al., 2024; Lucibello & Mézard, 2024). In the latter picture, each training sample contributes a local kernel, and generalisation arises from the superposition of many such kernels, see Figure 1a for a sketch. When kernels overlap strongly, the model assigns finite probability to the space between examples; when overlap is weak, generation can collapse onto individual training points. These analyses predict a *global* phase transition: the whole generative model goes from memorizing to generating new samples at once.

Recent work has begun to paint a more fine-grained picture of memorization. In practice, it was observed that models often reproduce only a subset of their training data rather than the entire dataset (Fang et al., 2025; Carlini et al., 2023); the effect depends strongly on the training setup (see further related work below). Thus, at finite sample size, a diffusion model often generates a mixture of copied and novel samples see Figure 1 c) f) and i). On the theoretical side, Achilli et al. (2026) showed that memorization may emerge progressively through the loss of manifold dimensions, leading to a form of *geometric memorization* in which some directions of variability are lost before exact copying occurs. Garnier-Brun et al. (2026) showed a coexistence of memorization and generalization in early-stopped diffusion models.

Taken together, these findings suggest that memorization is richer than a single global transition and raise a natural question:

**Which regions of data space, classes, or individual samples
are most likely to be memorized?**

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Figure 1. **Local data coverage governs memorization.** a) Sketch of local memorization in kernel density approximation. b) Histogram of data sparsity, measured by the average squared distance of a training point its 10 nearest neighbors, for 20,000 CIFAR10 samples. c) Generated samples + closest training image (measured by ℓ_2 distance) for a diffusion model trained on $P = 1000$ training examples. d) For each generated sample $\hat{x}$, we compute the local dominance as the log-ratio $\log d_2/d_1$ between the distances d_1 and d_2 to its nearest and second-nearest training samples. For each training sample x^μ, we show distances to generated images (red) and test data (blue), providing a baseline without memorization. Scatter shows raw values, markers indicate binned averages. e) Memorization rate per training sample x^μ, defined as the fraction of generated samples satisfying $d_1/d_2 < \tau$. Scatter shows raw values for $\tau = 0.1$, markers indicate binned averages across thresholds. f)–k) Same measurements for models trained on larger datasets. As the number of training samples increases, both dominance and memorization decrease, and their dependence on sparsity weakens. Overall, sparse regions exhibit strong single-sample dominance and higher memorization, while dense regions promote interpolation across multiple training points.

In this work, we give a concrete criterion for memorization by introducing a *local* theory of memorization in diffusion models. Our central hypothesis is simple: memorization is governed not only by the total number of training samples, but by their *local coverage*. Regions of data space containing many nearby examples support interpolation and generalization, whereas isolated regions are prone to sample retrieval. In Figure 1 we show an example of this behavior for diffusion models with U-net architecture (Ronneberger et al., 2015) trained on subsets of the CIFAR10 image dataset (Krizhevsky, 2009) of increasing size P. When trained on very few data points, the models first produce only memorized data. As we increase the number of training data, the models begin to generate a mixture of memorized and novel examples (panels c,f,i). We then quantify to which

degree samples generated from diffusion models replicate the closest training data (panels d,g,j) as a function of local sparsity. We find isolated training points in low density (high sparsity) regions are preferentially memorized (panels e,h,k), whereas samples generated in regions of higher local density (lower local sparsity in Figure 1) are less likely to be memorized. See Section 4 for full experimental details.

A local theory of memorization To formalize this intuition, we develop a local theory of memorization using the KDE framework of Lucibello & Mézard (2024); Biroli & Mézard (2024); Achilli et al. (2025b). Here, we sketch the main idea. We quantify the local coverage around a point $x \in \mathbb{R}^N$ by the probability mass contained in a ball of

radius h around x,

$$p_{\text{in}}(x) = \int_{B_h(x)} d\rho(x'), \qquad (1)$$

where $B_h(x)$ denotes a N-dimensional hypersphere centered at x and ρ the density of the data from which the training points are drawn. This quantity measures how densely the data distribution ρ populates the neighborhood of x. In the high-dimensional regime in which the number of samples scales as (Lucibello & Mézard, 2024; Biroli & Mézard, 2024)

$$P = e^{\alpha N}, \qquad (2)$$

we define

$$\nu_{\text{in}}(x) = \lim_{N \to \infty} \frac{1}{N} \ln p_{\text{in}}(x). \qquad (3)$$

Our main result shows that memorization at a point x is controlled by the pair $(\nu_{\text{in}}(x), \alpha)$ with a sharp transition:

$$\begin{cases} x \in A_{\text{mem}} & \text{if } \nu_{\text{in}}(x) < -\alpha \\ x \in A_{\text{gen}} & \text{otherwise} \end{cases} \qquad (4)$$

where we refer to A_{gen} as the region where the model correctly interpolates, and A_{mem} is a region where the learned density is characterized by isolated peaks centered on training examples. This result shows that depending on local coverage, the same diffusion model may exhibit retrieval-like behavior in some regions of space and generative interpolation in others. Points in low-coverage regions, where $\nu_{\text{in}}(x)$ is small, are memorized, while points in high-coverage regions are generalized.

Contributions In the following, we recall the relation between diffusion models and kernel density estimators in Section 2. Our main contributions are then as follows

- We develop **a local theory of memorization** and prove that the transition from memorization to generalization in a point is governed by the local kernel statistic ν_{in}, see Theorem 3.1 and Section 3.

- We show that the theory explains several empirical properties of diffusion models trained on CIFAR10 and CelebA (Liu et al., 2015) in Section 4:

  - **Coexistence of memorization and generalization:** the same model can simultaneously generate copied and novel samples; as P increases, memorization vanishes first in dense regions and persists longest in sparse regions, see Figure 1 and Section 4.1.
  - **Class-dependent memorization:** classes with larger intra-class diversity (and therefore lower local coverage) are memorized more strongly, see Figure 2 and Section 4.2.

More broadly, our results suggest that memorization in diffusion models is not a global binary property, but a local finite-sample phenomenon governed by the geometry and coverage of the training distribution. This mechanistic insight into memorization opens up the possibility of designing dedicated strategies to both suppress unwanted memorization and to retain specific training samples.

Further related Work

Impact of training setup on memorization A growing body of empirical and theoretical work has established that memorization depends strongly on properties of the training setup. In particular, it is exacerbated by data duplication (Carlini et al., 2023), prompt conditioning and classifier-free guidance (Somepalli et al., 2023; Wen et al., 2024; Kim et al., 2025; Gu et al., 2025), model capacity (Yoon et al., 2023; George et al., 2025; Zhang et al., 2025), dataset structure (Somepalli et al., 2023; Kim et al., 2025; Gu et al., 2025; Yoon et al., 2023), and training duration (Bonnaire et al., 2026; Favero et al., 2025; Garnier-Brun et al., 2026), while decreasing with learning rate (Wu et al., 2025) and dataset size (Somepalli et al., 2023; Kadkhodaie et al., 2024; Maillard & Goldt, 2026).

Local density estimators and memorization in practice Strategies to measure and mitigate memorization include determining the local intrinsic dimension of generated points (Ross et al., 2025), and estimating how "peaked" the estimated density is around a given point (Jeon et al., 2025) as well as identifying memorized prompts (Wen et al., 2024; Kim et al., 2025). These strategies share a common intuition: isolated data points produce a highly peaked local density, and produce samples of low intrinsic dimension. This perspective is naturally consistent with a kernel density estimation (KDE) view of diffusion models. Isolated training points induce highly concentrated local density estimates and generate low-dimensional samples, while dense regions support smoother interpolation. Similarly, conditioning can be interpreted as restricting the effective sample set contributing to the density estimate, thereby increasing local sparsity and promoting memorization. In this sense, existing empirical methods already implicitly rely on a local density perspective on memorization.

Memorisation in kernel density estimation On the theoretical side, several works have established a connection between diffusion models and kernel density estimation (Pidstrigach, 2022; Pham et al., 2024; Ambrogioni, 2024; Biroli & Mézard, 2024; Achilli et al., 2025a; Li et al., 2024; Lucibello & Mézard, 2024). Within this framework, one can determine a global "collapse phase" in diffusion models as a function of effective noise in the backward/sampling process (Biroli et al., 2024), where the model begins to

reproduce training data. More generally, the sample complexity required for smooth interpolation with kernel density estimators is known to scale exponentially with the ambient or manifold dimension(Lucibello & Mézard, 2024; Achilli et al., 2025b) of the data. Recent work has further shown that memorization can emerge gradually through a loss of manifold dimensions Achilli et al. (2026) and that this process may be spatially inhomogeneous. However, existing theoretical analyses remain largely global: they characterize when a model as a whole memorizes or generalizes, but do not predict which specific regions or samples are more likely to be memorized. In contrast, we develop a local theory of memorization that explicitly links memorization to data coverage at the level of individual points. This theory provides a principled explanation for the heterogeneous memorization patterns observed in practice.

2. Background: from diffusion models to kernel density estimation

We briefly recall a key observation: diffusion models trained on finite data behave like kernel density estimators (KDE). Diffusion models define a forward noising process that gradually corrupts a data point x_0 from ρ through the addition of noise

$$x_t\left(\epsilon_t\right) = \sqrt{\bar{\alpha}_t}x_0 + \sqrt{1 - \bar{\alpha}_t}\epsilon_t, \qquad \epsilon_t \sim \mathcal{N}\left(0, \mathrm{Id}\right), \quad (5)$$

where $\bar{\alpha}_t \in (0,1)$ is a decreasing function in t. As t increases, the original signal x_0 is gradually suppressed compared to the isotropic Gaussian noise, until one obtains x_T whose distribution is close to $\mathcal{N}\left(0, \mathrm{Id}\right)$. A neural network ϵ_θ with parameters θ is trained to predict the noise ϵ_t, which corresponds to learning the score $\nabla \ln \rho_t(x_t)$, where ρ_t is the distribution of x_t. This is achieved via the denoising score matching objective

$$L = \frac{1}{NTP}\sum_t\sum_{x^\mu \in \mathcal{D}}\mathbb{E}_{\epsilon_t}\left\|\epsilon_t - \epsilon_\theta\left(x_t^\mu\left(\epsilon_t\right), t\right)\right\|^2, \quad (6)$$

where $\mathcal{D}$ is a training set of size P points $\{x^\mu\}_{\mu=1}^P$ drawn i.i.d. from ρ.

To make the connection to KDE explicit, consider an idealized setting in which the model has infinite capacity and is trained to optimality. In this case, for each t, the learned function $\epsilon_\theta(\cdot, t)$ can be treated as an arbitrary function $\epsilon(\cdot, t)$ and the loss can be minimized functionally. Setting this functional derivative to zero yields:

$$\epsilon(x, t) \propto \nabla_x \ln \sum_\mu \mathcal{N}\left(x; \bar{\alpha}_t x^\mu, (1 - \bar{\alpha}_t)\mathrm{Id}\right). \quad (7)$$

This identity reveals that, when trained to convergence on a finite number of samples, diffusion models learn a Gaussian mixture centered at the training data points. In other words, diffusion models trained on a finite dataset recover the score of the empirical distribution convolved with Gaussian noise, rather than the true underlying density ρ. In the

low-noise limit ($\bar{\alpha}_t \to 1$), this mixture becomes sharply peaked around individual training samples, and the model effectively reproduces them. At finite noise, there is a direct correspondence between Equation (7) and KDE. KDE attempts to approximate a density ρ using a mixture of kernel functions K_N, indexed by the input dimension N, centered at the data points:

$$Z_N(x) = \frac{1}{P}\sum_\mu K_N\left(\frac{|x - x^\mu|}{h}\right). \quad (8)$$

The correspondence with KDE becomes explicit by identifying the kernel as Gaussian with bandwidth $h = \sqrt{1 - \bar{\alpha}_t}$ up to a rescaling of the variables x, x^μ by $\sqrt{\bar{\alpha}_t}$. In the ideal case, one has sufficient number of samples P such that Z becomes a smooth approximation of ρ. This behavior with P is one hypothesis to explain memorization and generalization in diffusion models.

This perspective suggests that generalization is not uniform across the data space. Even when the total number of samples is large, the quality of the KDE approximation depends on local sample density. Regions with high data density are well approximated, while sparse regions remain dominated by individual kernels. In the following, we formalize this intuition by deriving conditions under which local regions exhibit memorization or generalization.

3. Theory: Kernel density estimation with hard spheres

3.1. General theory and main result

To study kernel density estimation (KDE), we characterize the behavior of the log-density

$$z_N(x_N) = \frac{1}{N}\ln Z_N(x_N). \quad (9)$$

where $Z_N(x)$ is the mixture of kernel functions K centered at the data points defined in Equation (8). The difficulty in establishing the behavior of z_N is that it is a random quantity that depends on the draw of the P training points from ρ. Our goal is to compute its distribution over such draws to derive general properties of high dimensional KDE.

Since the ambient dimension changes with N, the symbol x_N should not be understood as a fixed point in a fixed space. We instead consider a sequence of vectors $\mathbf{x} = (x_N)_{N \geq 1}$ and distributions ρ_N with $x_N \in \mathbb{R}^N$, and assume that the corresponding local statistics have a limit. Equivalently, the result is indexed not by the coordinate vector x_N itself, but by the limiting local environment around it, namely the distribution of

$$\epsilon_{\mathbf{x}, N}(\tilde{x}_N) = \frac{1}{N}\ln K_N\left(\frac{x_N - \tilde{x}_N}{h}\right), \qquad \tilde{x}_N \sim \rho_N,$$

Thus, the behavior of $z_N(x_N)$ is controlled by the distribution of the scalar statistic $\epsilon_{\mathbf{x},N}$ rather than by the full high-dimensional variable $\tilde{x}_N \sim \rho_N$.

We use an exponential transform, inspired by Dotsenko (2011), to compute the cumulative distribution function of $z_N(x_N)$. Define

$$\overline{W_N(y, x_N)} = \overline{\exp\left(-\exp\left[-N\left(y - z_N(x_N)\right)\right]\right)}, \quad (10)$$

where the overline denotes the average over draws of the P training samples. As $N \to \infty$, the inner function converges pointwise to $e^{-e^{-N(\cdot)}} \to \Theta(\cdot)$, with convention $\Theta(0) = e^{-1}$. Hence if the limiting CDF $F_{\mathbf{x}}(y) = \lim_{N\to\infty} \Pr(z_N(x_N) \leq y)$ is continuous in y, then $\overline{W_N(y, x_N)}$ converges to F as $N \to \infty$; see Section A.1.

Using independence of the training samples, the average factorizes and gives

$$\overline{W_N(y, x_N)} = \exp\left\{e^{\alpha N} \ln\left[1 - I_N(y, x_N)\right]\right\}, \quad (11)$$

with

$$I_N(y, x_N) = \mathbb{E}_{\tilde{x}\sim\rho_N}\left[1 - e^{-e^{N\left[\epsilon_{x,N}(\tilde{x}) - \alpha - y\right]}}\right]. \quad (12)$$

This yields the following theorem, which is our main theoretical result:

Theorem 3.1 (Cumulative distribution function of KDE). *Assume that the cumulative distribution function converges $F_{\mathbf{x},N}(y) = \Pr(z_N(x_N) \leq y) \to F_{\mathbf{x}}(y)$ for all y as $N \to \infty$. Furthermore, let us assume that the limit*

$$i_{\mathbf{x}}(y) := \lim_{N\to\infty} e^{\alpha N} I_N(y, x_N) \quad (13)$$

exists with $i_{\mathbf{x}}(y) \in [0, \infty]$. Then if $F_{\mathbf{x}}(y)$ is continuous in y, we find

$$F_{\mathbf{x}}(y) = \begin{cases} 0 & \text{if } i_{\mathbf{x}}(y) = \infty, \\ \exp[-i_{\mathbf{x}}(y)] & \text{if } 0 < i_{\mathbf{x}}(y) < \infty, \\ 1 & \text{if } i_{\mathbf{x}}(y) = 0. \end{cases} \quad (14)$$

We detail the proof in Section A.

Theorem 3.1 shows that the limiting CDF of $z_N(x_N)$ is determined by the exponential scaling of the probability that the kernel statistic $\epsilon_{\mathbf{x},N}$ is large enough to contribute at level y. In typical high-dimensional settings, $z_N(x_N)$ concentrates, so the intermediate regime in Equation (14) collapses to the location of a sharp jump. Fluctuations of $\epsilon_{\mathbf{x}}$ would instead broaden this transition.

Relation to Gumbel's law. The form of Equation (11) is somewhat reminiscent of Gumbel's law which computes the cumulative distribution function of the maximum of

P i.i.d. random variables, in our case, this amounts to replacing $z(x)$ by $\tilde{z}(x) = \ln \max_\mu K\left(\frac{x-x^\mu}{h}\right)/P$. This structural similarity is no accident, indeed one can show that whenever $z(x) \approx \tilde{z}(x)$, Equation (11) directly corresponds to Gumbel's law. In the generalization regime however, exponentially many terms contribute to z, and the analogy breaks down. We discuss this relation in more detail in Section A.4.

3.2. Case study: hard-sphere kernels

We now specialize to the hard-sphere kernel

$$K_N(u) = \frac{\Theta(R^2 - \|u\|^2)}{V_R}, \quad (15)$$

where V_R is the volume of the N-dimensional ball of radius R. In high dimensions, the hard-sphere kernel captures the same radial concentration mechanism as a Gaussian kernel. At the level of the exponential transform, the relevant quantity is the distribution of $\epsilon_{x,N}(\tilde{x}_N)$. For a Gaussian kernel, this statistic is, up to additive constants, proportional to $-\frac{1}{N}\|x_N - \tilde{x}_N\|^2$. Thus the leading exponential behavior is governed by level sets of the squared distance. In high dimensions, the Gaussian measure concentrates on a thin shell at its typical radius. Similarly, for the hard-sphere kernel, almost all of the volume of the ball concentrates near its boundary. Therefore, while Gaussian and hard-sphere kernels are quantitatively different, the hard-sphere kernel provides a tractable model that exposes the same local-coverage transition mechanism. For simplicity, we fix $R = 2\pi e$, for which the volume V_R remains $\mathcal{O}(1)$ as $N \to \infty$. The effective scale of the kernel relative to the data is controlled by the bandwidth h, which can be interpreted as a rescaling of the data points by a factor $1/h$. Unless otherwise specified, we set $h = 1$.

In this case, we find that the quantity controlling the limiting behavior is the local coverage. Let us assume that the limit of the local coverage rate

$$\nu_{\text{in},\mathbf{x}} = \lim_{N\to\infty} \frac{1}{N} \ln p_{\text{in},N}(x_N). \quad (16)$$

exists. Then we obtain the following result:

Lemma 3.2 (Local hard-sphere KDE transition). *Let $P = e^{\alpha N}$ with fixed $\alpha > 0$, and let K_N be the hard-sphere kernel in Equation (15). Under the assumptions of Theorem 3.1 and additionally assuming that that $\nu_{\text{in},\mathbf{x}}$ exists, when $\nu_{\text{in},\mathbf{x}} \neq -\alpha$ we find*

$$z_N(x_N) \xrightarrow[N\to\infty]{\mathbb{P}} \begin{cases} \nu_{\text{in},\mathbf{x}} & \text{if } \nu_{\text{in},\mathbf{x}} > -\alpha, \\ -\infty & \text{if } \nu_{\text{in},\mathbf{x}} < -\alpha. \end{cases} \quad (17)$$

The proof follows from a direct application of Theorem 3.1, see Section A.

Lemma 3.2 establishes a local transition controlled by the comparison between local coverage $\nu_{\text{in}}(x)$ and sample complexity α. Regions with $\nu_{\text{in}}(x) > -\alpha$ contain exponentially many nearby samples and therefore support smooth KDE estimates. Regions with $\nu_{\text{in}}(x) < -\alpha$ contain no nearby samples with high probability; in the diffusion-model interpretation, probability mass assigned to such regions concentrates on isolated training examples, leading to memorization. The divergence $z(x) \to -\infty$ for $\nu_{\text{in}}(x) < -\alpha$ reflects the fact that, in these regions, the kernel density estimate vanishes with high probability. Indeed,

$$\Pr\left(\{Z(x) = 0\}\right) = (1 - p_{\text{in}}(x))^P$$
$$= e^{P \ln(1 - e^{N\nu_{\text{in},N}(x)})}$$
$$\to \Theta\left(-[\alpha + \nu_{\text{in},\mathbf{x}}]\right)$$

On the other hand, when α is large enough, then $Z(x)$ approaches $p_{\text{in}}(x)$, hence the kernel density estimate converges to a local average of ρ over a sphere.

In the context of diffusion models, such regions are dominated by isolated training points. If probability mass is assigned to these locations, it must concentrate on individual samples, leading to memorization. This establishes the coexistence of memorized and generalized regions, as described in Eq. (4).

4. Experimental results

4.1. Local sparsity governs memorization of samples

We now empirically investigate how local data density influences memorization behavior in diffusion models. Our goal is to test the central prediction of the KDE-based theory that sparse regions are more prone to memorization than dense ones. This implies a coexistence of memorization and generalization: copied and novel samples can arise simultaneously from the same model. Furthermore, it predicts a gradual, rather than abrupt disappearance of memorization: as P increases, memorization vanishes first in dense regions and persists longest in sparse regions. Consequently, regions of low data density lead to more localized, sample-specific behavior, while dense regions promote interpolation across multiple training points. While our theory predicts a sharp transition as a function of local coverage, here we test its qualitative consequences using empirical proxies for density.

4.1.1. EXPERIMENTAL METHODS

We train diffusion models with a U-net architecture on randomly chosen subsets of CIFAR-10 and CelebA. We then generate samples from these models and assess their closeness to training examples using different measures outlined below. Details on the experimental procedure are given in

Section C. We first define how we measure sparsity and memorization and then describe the outcome of the experiments shown in Figure 1 and Figure 3.

Local sparsity. For each training sample x^μ, we quantify local sparsity using nearest-neighbor distances, which serve as a proxy for inverse local coverage. Concretely, we define

$$s(x^\mu) = \frac{1}{Nk} \sum_{i=1}^{k} \|x^\mu - x^{\mu_i}\|^2, \tag{18}$$

where $\{x^{\mu_i}\}$ are the k nearest neighbors of x^μ. Larger values of $s(x^\mu)$ correspond to sparser regions of the data distribution. We also validate that the results remain consistent for different choices of k in $\{5, 10, 20, 50\}$. The results reported in the figures correspond to $k = 10$.

Local dominance. Given a generated sample $\hat{x}$, we compute its distances $d_1(\hat{x})$ and $d_2(\hat{x})$ to the nearest and second-nearest training samples, and define the *dominance*

$$\Delta(\hat{x}) = \log \frac{d_2(\hat{x})}{d_1(\hat{x})}. \tag{19}$$

Large values of Δ indicate that a single training point dominates the local score estimation around the generated sample, while $\Delta \approx 0$ corresponds to interpolation between multiple neighbors. We assign each generated sample to its nearest training point and compute, for each x^μ, the average dominance over all generated samples assigned to it. To disentangle model-specific effects from dataset geometry, we construct a baseline by replacing generated samples with held-out test data, processed in the same way.

Memorization rate. We further report a binary memorization metric used in several previous studies (Wu et al., 2025; Yoon et al., 2023; Bonnaire et al., 2026) based on the condition $d_1/d_2 < \tau$.

4.1.2. RESULTS

The middle column of Figure 1 shows the average dominance as a function of local sparsity for different training set sizes. We observe a clear **monotonic increase of dominance with sparsity** for generated samples: sparse regions exhibit strong single-sample dominance, while dense regions show low dominance and interpolation. In contrast, the test-data baseline remains nearly flat, indicating that this effect is not explained by nearest-neighbor geometry alone. As the number of training samples increases, the overall level of dominance decreases and the dependence on sparsity weakens, consistent with the expectation that higher sample density reduces isolated regions. The corresponding results for the memorization rate are shown in the right column of Figure 1. While we do not directly estimate the threshold $\nu_{\text{in}}(x) = -\alpha$, the observed dependence on local sparsity is consistent with a transition controlled

by local coverage. For the canonical value $\tau = 0.33$ commonly used in the literature, we see a clear **coexistence of memorization and generalization** that depends on data sparsity. Moreover, while the precise behavior depends on τ, the same qualitative trend is observed: **memorization increases with sparsity**. In the appendix Section B we show the outcome of the same experiment on downsampled CelebA, an image dataset of celebrity portraits. The results of these experiments are also consistent with the hypothesis. These results support a local version of the KDE picture: diffusion models exhibit a continuous transition from interpolation in dense regions to single-sample dominance in sparse regions.

4.2. Intra-class sparsity predicts class-wise memorization

We now study memorization at a coarser level, focusing on **class-dependent effects**. In a multi-class setting, the KDE picture predicts that classes with more concentrated support (i.e. lower local sparsity) should be memorized less than classes with higher intra-class variability. A schematic illustration is shown in Figure 2a. To formalize this intuition, consider a mixture of class-conditional densities ρ_c, such that

$$\rho(x) = \sum_c w_c \rho_c(x),$$

where the weights w_c sum to one, and we assume that all weights are order one in N. We define the local in-distribution log-density for class c as

$$\nu_{\text{in},c}(x) := \frac{1}{N} \ln p_{\text{in},c}(x) = \frac{1}{N} \ln \int_{B_h(x)} d\rho_c(x'), \quad (20)$$

and assume that $\nu_{\text{in},c}(x)$ remains $\mathcal{O}(1)$ in the large-N limit. Then the total in-distribution density satisfies

$$\nu_{\text{in}}(x) = \frac{1}{N} \ln \sum_c e^{N\left(\nu_{\text{in},c}(x) + N^{-1} \ln w_c\right)} \longrightarrow \max_c \nu_{\text{in},c}(x),$$

i.e. it is dominated by the locally densest class. Consequently, we obtain the same scaling behavior as in lemma 3.2, in a class-dependent manner:

$$z(x) \rightarrow \begin{cases} \max_c \nu_{\text{in},c}(x) & \text{if } \max_c \nu_{\text{in},c}(x) \geq -\alpha, \\ -\infty & \text{otherwise}, \end{cases}$$
$$(21)$$

showing that memorization is governed by the class with the highest local coverage around x. This leads to a clear prediction: classes with higher intra-class variability (and thus lower local coverage) should be more prone to memorization.

We now demonstrate empirically that such class-dependent density differences naturally arise. As a concrete example, we isolate the memorization behavior of different subclasses of CIFAR-10. In Figure 2b we show that the classes

"airplane" and "ship" in CIFAR-10 have, on average, considerably lower intra-class diversity than the classes "truck" and "automobile", likely due to more homogeneous blue background colors in the former two classes. Correspondingly, we observe both higher local dominance, and higher memorization in more diverse classes.

In the appendix, Section B, we show that the same trends emerge for two additional experiments. First, when one sorts the CelebA datasets into classes according to whether images have the attribute "wearing hat" or "blond", images where "wearing hat" is true are more diverse than those where it is false. Correspondingly, portraits featuring hats are memorized more.

Second, we train diffusion models on a mixture of MNIST (Deng, 2012) and CIFAR-10 images. MNIST, consisting of handwritten digits, is structurally simpler and occupies a much more concentrated region of image space, whereas CIFAR-10 exhibits substantially higher variability. Treating MNIST and CIFAR-10 as two distinct classes, we observe that MNIST samples consistently exhibit lower local sparsity than CIFAR-10 samples. As predicted by the theory, this translates into reduced memorization. These results support the hypothesis that classes with higher intra-class diversity are memorized more strongly than classes with lower intra-class diversity.

5. Discussion

Summary. In this work, we test the predictive power of kernel density estimation to explain local memorization in diffusion models. While earlier works (Lucibello & Mézard, 2024; Biroli & Mézard, 2024; Achilli et al., 2025b; 2026) focused on establishing *global* phase transitions from memorization to generalization for *typical* samples from the data distribution, we developed a *local* theory of memorisation for KDE. We showed that, in the high-dimensional limit, each point x in sample space undergoes a local transition from memorisation to generalisation as a function of sample complexity. This transition is governed by *local coverage* of training data, which separates regions where the density is smoothly estimated from regions dominated by isolated training samples. This leads to a coexistence of *generalization* and *memorization* within the same model. Our analysis further predicts that, in multi-class settings, the behavior at a point x is controlled by the class with the highest local coverage. Empirically, we find consistent evidence for these predictions: memorization increases with local sparsity and intra-class variability, and different classes exhibit systematically different memorization behavior. Taken together, these results show that memorization in diffusion models is not a global failure mode, but a *localized phenomenon* driven by the geometry of the data distribution. In particular, isolated regions of the data space can remain memorized

Figure 2. **Intra-class sparsity predicts class-wise memorization.** a) Schematic illustration of KDE for two classes with different local sparsities. b) Distribution of per-class sparsity around training points for different subclasses of CIFAR-10, measured from 20,000 training samples. c) Generated samples together with their closest training example (measured by cosine similarity) for a diffusion model trained on $P = 1000$ samples, sorted to contain 2 nearest neighbors for each of the four classes shown in b). d) Local dominance as a function of sparsity. For each generated sample $\hat{x}$, we compute distances d_1 and d_2 to its nearest and second-nearest training samples, and define dominance via their ratio. Each point corresponds to a training sample x^μ, aggregating generated samples assigned to it. Scatter shows raw values, markers indicate binned averages. Different colors correspond to different classes. e) Memorization rate of each sample x^μ, defined as the fraction of generated samples satisfying $d_1/d_2 < \tau = 1/3$. f) - k) report the same measures, but for diffusion models trained on larger datasets. "Airplane" and "ship" samples lie in denser regions and exhibit lower dominance and memorization, while "truck" and "automobile" samples are sparser and more frequently memorized, consistent with the prediction that local coverage controls class-dependent memorization.

even when the majority of generated samples are novel.

Limitations. Our analysis relies on the infinite-dimensional limit $N \to \infty$ with an exponential sample complexity $P = e^{\alpha N}$, analogous to prior work (Lucibello & Mézard, 2024; Biroli & Mézard, 2024). While this enables a sharp theoretical characterization, it predicts an abrupt transition between memorization and generalization as a function of local coverage, whereas empirically we observe a more gradual dependence on local sparsity. This more gradual behavior could be the result of finite-size fluctuations as well as an extended region where the second condition in Theorem 3.1 is fulfilled, which allows for fluctuations even in the infinite limit. While we hypothesize that finite-N fluctuations are at the core of this more gradual

increase, disentangling these effects from fluctuations which persist in the infinite limit is an important direction for future work. Similarly, the qualitative equivalence between hard-sphere kernels and Gaussian kernels, which we exploit here to make the dependence of KDE on local coverage explicit, is expected to increase with the dimension N. We leave an investigation into the effects of the exact shape of the kernel to future work.

A second limitation stems from the choice of the metric space underlying the KDE approximation. Our theory assumes that diffusion models operate in the ambient space. In practice, however, models may implicitly operate in a lower-dimensional representation, for example through projection onto a data manifold (Achilli et al., 2025a; 2026) or by exploiting symmetries in the data (Kamb & Ganguli,

2025), or as latent diffusion models (Rombach et al., 2022). In such cases, the effective kernel should be understood as acting in this learned metric space, which depends on model architecture and may help explain the observed dependence of memorization on model capacity (Yoon et al., 2023), which has already been shown explicitly for autoencoder architectures by Zhang et al. (2025).

Outlook. Our results suggest that controlling memorization requires shaping the *local geometry* of the data representation. In particular, learning representations that increase local coverage, by mapping data to lower-dimensional spaces or by exploiting invariance may reduce memorization but reduces model expressivity, presenting a tradeoff between increased local coverage and preserving generative performance.

Another important direction concerns training dynamics. Memorization is known to depend on training time and optimization hyperparameters (Bonnaire et al., 2026; Favero et al., 2025; Wu et al., 2025), and recent work suggests that memorization and generalization can emerge simultaneously during training (Garnier-Brun et al., 2026). Understanding how convergence to the KDE solution in the generalized and memorized regimes depends on optimization dynamics is a promising direction for future research.

Broader impact. While this work is foundational, memorization in generative models is directly related to privacy and copyright risks. Understanding which examples are likely to be memorized may help audit and mitigate training-data reproduction, although the same insights could potentially inform extraction attempts. At the same time, a better understanding of memorisation might help devise strategies to deliberately store samples in diffusion model.

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A. Details on Theoretical results

In this appendix, we give a detailed proofs of the results presented in Section 3.

A.1. Convergence of the exponential transform

To establish the behavior of the KDE, we characterize the cumulative distribution function (CDF) of the log-density z through the exponential transform Equation (10). At the base of Theorem 3.1 therefore lies the convergence of the exponential transform to the CDF, which is provided by the following Lemma.

Lemma A.1 (Convergence of the exponential transform to the CDF.). *Let $N \in \mathbb{N}, N > 0, x_N \in R^N$ and $F_N(y) = \Pr(z_N(x_N) \le y)$ be the cumulative distribution function of $z_N(x_N)$. Then for $\overline{W_N(y, x_N)}$ defined as in Equation (10) and $\delta > 0$, we find that*

$$e^{-e^{-N\delta}} F_N(y - \delta) \le \overline{W_N(y, x_N)} \le F_N(y + \delta) + e^{-e^{N\delta}} \Pr(z_N \ge y + \delta).. \tag{22}$$

Consequently, if the limit $F_N \to F_{\mathbf{x}}$ exists and if $F_{\mathbf{x}}$ is continuous in y, then

$$\overline{W_N(y, x_N)} \to F_{\mathbf{x}}(y).$$

Proof. For notational simplicity we suppress the dependence on x and write

$$z_N := z_N(x), \qquad W_N(y) := \exp\{-\exp[-N(y - z_N)]\}.$$

We first show the lower bound. If $z_N \le y - \delta$, then

$$W_N(y) = \exp\{-\exp[-N(y - z_N)]\} \ge \exp\{-e^{-N\delta}\}.$$

In other words, using the indicator function $\mathbf{1}$, we know that

$$W_N(y) \ge e^{-e^{-N\delta}} \mathbf{1}\{z_N \le y - \delta\}.$$

Taking expectations gives

$$\overline{W_N(y)} \ge e^{-e^{-N\delta}} \mathbb{P}(z_N \le y - \delta).$$

We now take care of the upper bound. Since $0 \le W_N(y) \le 1$, we know that

$$W_N(y) \le \mathbf{1}\{z_N < y + \delta\} + W_N(y)\mathbf{1}\{z_N \ge y + \delta\}$$
$$\le \mathbf{1}\{z_N < y + \delta\} + e^{-e^{N\delta}} \mathbf{1}\{z_N \ge y + \delta\}.$$

Taking the average yields

$$\overline{W_N(y)} \le \mathbb{P}(z_N < y + \delta) + e^{-e^{N\delta}} \mathbb{P}(z_N \ge y + \delta).$$

Combining the two bounds yields Lemma A.1. □

A.2. Average over draws of samples Theorem 3.1

We now proceed with the proof of Theorem 3.1. Let us first keep $N \in \mathbb{N}, N > 0, x_N \in R^N$ fixed. Our starting point is Equation (10). Recall that the notation $\overline{f(z)}$ denotes the average over the data set, meaning that when we insert the definition of $z(x)$ and explicitly state the average, we find

$$\overline{W_N(y, x_N)} = \int \prod_\mu d\rho(x^\mu) \exp\left(-e^{-N(y+\alpha)} \sum_\mu K(x - x^\mu)\right)$$

$$= \left\{\int d\rho(\tilde{x}) \exp\left[-e^{-N(y+\alpha)} K(x - \tilde{x})\right]\right\}^P$$

which directly yields Equation (11). Taking the limit then reveals the three scaling regimes

$$\overline{W_N(y, x_N)} \to \begin{cases} 0 & \text{if } i_{\mathbf{x}}(y) = \infty, \\ \exp[-i_{\mathbf{x}}(y)] & \text{if } 0 < i_{\mathbf{x}}(y) < \infty, \\ 1 & \text{if } i_{\mathbf{x}}(y) = 0. \end{cases} \tag{23}$$

Using, lemma A.1, we then find that the limit of the exponential transform yields the limiting CDF. □

A.3. Hard spherical kernels

We now turn to the proof of lemma 3.2. Using the definition of p_{in}, we find that the integral I_N defined in Equation (12) decomposes into two areas: those where $K = 0$ and those where $K = 1/V_R$. Using this distinction, we then find

$$I_N(y, x_N) = p_{\text{in},N}(x_N)\left(1 - \exp\left(-\exp\left\{-N\left(y + \alpha + \frac{1}{N}\ln V_R\right)\right\}\right)\right) \tag{24}$$

Since, we have fixed R such that $\frac{1}{N}\ln V_R \to 0$, the latter term in the exponential vanishes in the high-dimensional limit. To apply Theorem 3.1, we use the assumption that $\nu_{\text{in},\mathbf{x}}$ exists and compute the limit Equation (13) in two distinct regimes: $y + \alpha > 0$ and $y + \alpha < 0$.

In the first regime, $y + \alpha = \delta > 0$, we find

$$i_{\mathbf{x}} = \lim_{N \to \infty} e^{N(\alpha + \nu_{\text{in},N}(x_N))}\left(1 - e^{-e^{-N\delta}}\right) = \begin{cases} 0 & \text{if } y - \nu_{\text{in},\mathbf{x}} > 0 \\ \infty & \text{if } y - \nu_{\text{in},\mathbf{x}} < 0 \end{cases}.$$

In the second regime, $y + \alpha = -\delta < 0$, we find

$$i_{\mathbf{x}} = \lim_{N \to \infty} e^{N(\alpha + \nu_{\text{in},N}(x_N))}\left(1 - e^{-e^{N\delta}}\right) = \begin{cases} 0 & \text{if } \alpha + \nu_{\text{in},\mathbf{x}} < 0 \\ \infty & \text{if } \alpha + \nu_{\text{in},\mathbf{x}} > 0 \end{cases}.$$

Applying Theorem 3.1 to both regimes, we find

$$F_{\mathbf{x}}(y) = \begin{cases} \Theta(y - \nu_{\text{in},\mathbf{x}}) & y + \alpha > 0 \\ \Theta(-(\nu_{\text{in},\mathbf{x}} + \alpha)) & y + \alpha < 0 \end{cases}$$

which yields the probability that $z \leq y$ in the limit, thanks to lemma A.1. The first line shows that when y is larger than $-\alpha$, the CDF becomes a step function with a jump at $y = \nu_{\text{in},\mathbf{x}}$.

We split the second line into two scenarios, $\nu_{\text{in},\mathbf{x}} + \alpha > 0$ and $\nu_{\text{in},\mathbf{x}} + \alpha < 0$ In the first, scenario $\nu_{\text{in},\mathbf{x}} > -\alpha > y$, hence this implies $y < \nu_{\text{in}}(x)$, and the limiting CDF is zero in y. On then other hand, in the second scenario, when $\nu_{\text{in},\mathbf{x}} < -\alpha$, then the CDF is one for any finite value of y. This means that z is smaller than any finite value of y with probability one. Taken together, the first and the second line imply that the CDF has a jump at $y = \nu_{\text{in},\mathbf{x}}$ when $\nu_{\text{in},\mathbf{x}} > -\alpha$. The second scenario in the second line implies that when $\nu_{\text{in},\mathbf{x}} < -\alpha$, all draws of z are smaller than any finite y, hence $z \to -\infty$.

Consequently, we find that the cumulative distribution has the shape of a step function. Hence the value of $z_N(x_N)$ concentrates on the "jump location", given by

$$\begin{cases} \nu_{\text{in}}(x) & \text{if } \nu_{\text{in}}(x) \geq -\alpha \\ -\infty & \text{if } \nu_{\text{in}}(x) < -\alpha \end{cases} \tag{25}$$

which finally yields lemma 3.2. $\qquad\square$

A.4. Relation to Gumbel-type extreme value statistics

In this appendix we clarify the relation between the exponential transform used in Eq. (10) and Gumbel-type extreme value statistics. Let us first clarify what we mean by Gumbel-type extreme value statistics: for independent random variables $u_1, \ldots, u_P$, one has

$$\mathbb{P}\left(\max_{\mu} u_{\mu} < y\right) = [1 - \mathbb{P}(u_{\mu} \geq y)]^P. \tag{26}$$

which is the standard structure associated with Gumbel-type extreme value statistics and, identifying $I_{\circ}$ with $\mathbb{P}(u_{\mu} \geq y)$, has strong similarity with Equation (11).

We will first show that letting $N \to \infty$ at $P < \infty$ fixed, W_N has a natural extreme-value interpretation. However, this does not mean that the calculation replaces the KDE by its largest term, because in the main result, we let $P \to \infty$ with $\frac{1}{N}\ln P = \alpha$ fixed. In this case, W_N is a soft probe of the cumulative distribution function of the full log-KDE.

For the sake of comparison, however, let us now keep $P < \infty$ and all samples fixed. We then find that W_N can be rewritten as

$$W_N(y, x) = \prod_{\mu=1}^{P} \exp\left\{-K(x - x^\mu)e^{-N(y+\alpha)}\right\}. \tag{27}$$

Now define the single-sample contribution $u_\mu(x) := \frac{1}{N} \ln K(x - x^\mu) - \alpha$. Then each factor in Eq. (27) can be written as

$$\exp\left\{-K_\mu(x)e^{-N(y+\alpha)}\right\} = \exp\left\{-\exp[N(u_\mu(x) - y)]\right\}. \tag{28}$$

If we were to fix P and all $u_\mu(x)$, and let $N \to \infty$ this converges to a hard threshold:

$$\exp\left\{-\exp[N(u_\mu(x) - y)]\right\} \xrightarrow[N\to\infty]{} \Theta(y - u_\mu(x)). \tag{29}$$

Consequently, if this hard-threshold limit is taken termwise before accounting for the exponentially many samples, Eq. (27) becomes

$$W_N(y, x) \approx \prod_{\mu=1}^{P} \Theta(y - u_\mu(x)) = \Theta\left(y - \max_\mu u_\mu(x)\right). \tag{30}$$

This is the origin of the formal similarity with Gumbel's law: the product form resembles the cumulative distribution function of a maximum. In regimes where the KDE is dominated by a single large kernel contribution, $z_N(x)$ is well approximated by

$$\widetilde{z}_N(x) := \max_\mu \frac{1}{N} \log \frac{K_\mu(x)}{P} = \max_\mu u_\mu(x). \tag{31}$$

In such max-dominated regimes, the above Gumbel-type interpretation directly describes the statistics of $z_N(x)$.

However, to obtain a generalization regime, we must send P to infinity too. Consequently, in the generalization regime, exponentially many samples may contribute to the KDE at any point. Then the sum cannot be replaced by its largest term. This distinction can be seen explicitly for the hard-sphere kernel. Let $m_N(x)$ be the number of training samples inside the corresponding ball around x with volume $V_R = 1$. Then

$$Z_N(x) = \frac{m_N(x)}{P}, \qquad \Rightarrow \qquad z_N(x) = \frac{1}{N} \log m_N(x) - \alpha. \tag{32}$$

If the local mass of the ball satisfies $\ln p_{\text{in}}(x) = N\nu_{\text{in}}(x) + o(N)$, then the typical number of samples in the ball is

$$m_N(x) \sim Pp_{\text{in}}(x) = e^{N(\alpha + \nu_{\text{in}}(x)) + o(N)} \tag{33}$$

whenever $\alpha + \nu_{\text{in}}(x) > 0$. In this regime,

$$z_N(x) \to \nu_{\text{in}}(x), \tag{34}$$

up to subexponential prefactors. By contrast, the largest single contribution is only

$$\widetilde{z}_N(x) = -\alpha \tag{35}$$

provided that at least one sample lies inside the ball. Thus, whenever $\nu_{\text{in}}(x) > -\alpha$, the full KDE exponent $z_N(x)$ differs from the maximum contribution $\widetilde{z}_N(x)$.

The Gumbel-like structure of Eq. (11) should therefore be understood as follows: the exponential transform is a soft thresholding device which, after factorization over samples, resembles an extreme-value observable. If one hardens this threshold before accounting for the exponentially many samples, one obtains a maximum-probe interpretation. Keeping the transform soft through the large-N calculation, however, retains the collective contribution of exponentially many moderate terms and therefore probes the full log-KDE $z_N(x)$.

B. Additional Experiments

B.1. Local sparsity and dominance

In Figure 3, we report the analogous experiment to Figure 1 for CelebA data, which we downsample to 32×32 greyscale pixels. We find that both local dominance and memorization increase with local sparsity. In comparison to CIFAR-10, we observe that the CelebA dataset appears to have fewer samples in very low density regions (compare logarithmic scale of Figure 3 d) to Figure 1d).

Figure 3. **Local memorization in CelebA data.** a) Sketch of kernel density approximation and local memorization phenomenon. b) Distribution of data sparsity for 40,000 samples from CelebA. c) Generated samples + closest training image (measured by cosine similarity) for diffusion model trained on $P = 4000$ training examples. d) Local dominance as a function of sparsity. For each generated sample $\hat{x}$, we compute distances d_1 and d_2 to its nearest and second-nearest training samples, and define dominance via their ratio. Each point corresponds to a training sample x^μ, aggregating generated samples assigned to it. Scatter shows raw values, markers indicate binned averages. Blue points show the same quantity computed using test data, providing a baseline without memorization. e) Memorization rate per training sample x^μ, defined as the fraction of generated samples satisfying $d_1/d_2 < \tau$. Scatter shows raw values for $\tau = 0.1$, markers indicate binned averages across thresholds. f)–k) Same measurements for models trained on larger datasets. As the number of training samples increases, both dominance and memorization decrease, and their dependence on sparsity weakens. Overall, sparse regions exhibit strong single-sample dominance and higher memorization, while dense regions promote interpolation across multiple training points.

B.2. Attribute-dependent Memorization, CelebA

Figure 4. **Attribute-conditioned memorization in CelebA.** a) Generated samples + closest training image (measured by cosine similarity) for diffusion model trained on $P = 20000$ training examples, sorted by classes. b) Distribution of data sparsity for 40000 samples from CelebA. d) Local dominance of training sample x^μ against local data sparsity around x^μ for a diffusion model trained on $P = 20000$ training examples from CelebA.

The CelebA dataset consists of celebrity portraits that are annotated with "attributes" such as hair color or accessories. We construct four classes from these attributes, conditioning on "blond" and "not blond" as well as "wearing hat" and the opposite. Again, we find that local dominance (and thus memorization) increases with data sparsity, see Figure 4. Samples from classes where "wearing hat" is true are typically more diverse. Therefore this class has a higher average sparsity per class,and is therefore more likely to be memorized. These classes are not balanced: approximately one sixth of images has attribute "blond" and approximately 5 % of images have attribute "wearing hat". Consequently the diffusion model generates fewer samples that are memorized data points with these attributes, and even fewer samples are closest to training data where both attributes are true. This is reflected in the lower bin resolution in Figure 4b) and fewer such points appearing in Figure 4 c).

B.3. Class-dependent Memorization, CIFAR-10+MNIST

In a separate experiment, we train diffusion models on a mixture of MNIST (Deng, 2012) and CIFAR-10 images. MNIST, consisting of handwritten digits, is structurally simpler and occupies a much more concentrated region of image space, whereas CIFAR-10 exhibits substantially higher variability. Treating MNIST and CIFAR-10 as two distinct classes, we observe that MNIST samples consistently exhibit lower local sparsity than CIFAR-10 samples. As predicted by the theory, this translates into reduced memorization: in Figure 5, both local dominance and memorization rates increase with sparsity. For the comparison of CIFAR and MNIST, all distances are computed on ℓ_2-normalized samples, so that sparsity reflects relative geometric structure (angular similarity) rather than differences in overall scale. We verified that using unnormalized samples yields qualitatively similar results. On average, MNIST points lie in denser regions and are therefore memorized less than CIFAR-10 points.

C. Experimental Details

Model. We train denoising diffusion models using a standard discrete-time formulation with $T = 1000$ timesteps. The score network is parameterized by a U-Net architecture (Ronneberger et al., 2015) with convolutional filters. All models operate directly in pixel space.

Training procedure. Models are trained using the standard diffusion objective with mean squared error loss. Optimization is performed using Adam with learning rate 10^{-4} and batch size 100. Training proceeds for a fixed number of 10^6 gradient steps for all models.

Figure 5. **Class-wise memorization on combined MNIST and CIFAR-10.** a) Generated samples together with their closest training example (measured by cosine similarity) for a diffusion model trained on $P = 2000$ samples. b) Distribution of per-class sparsity around training points for four classes in CIFAR-10 measured from 20,000 training samples. c) Generated samples together with their closest training example (measured by cosine similarity) for a diffusion model trained on $P = 5000$ samples. d) Class-wise local dominance of a training point x^μ as a function of local data sparsity. Local dominance is defined via the ratio between the distance d_1 to the closest generated sample and the distance d_2 between that generated sample and the second-closest training point. MNIST points are shown in blue, CIFAR-10 points in red. e) Memorization rate of each sample x^μ, defined as the fraction of generated samples satisfying $d_1/d_2 < \tau = 1/3$. Different classes are shown as different colors. f)–k) Same measurements for models trained on larger datasets. As the number of training samples increases, both dominance and memorization decrease, and their dependence on sparsity weakens.

Evaluation   For all experiments, the dataset is deterministically split into a training set of size P and a held-out test set. The same split is reused for all evaluations to ensure comparability across checkpoints and metrics. At evaluation time, we generate $N_{\text{gen}} = 4000$ samples from the trained diffusion model. We compare generated samples to training and test data using either cosine similarity or normalized ℓ_2 distance. In the cosine case, all inputs are flattened and ℓ_2-normalized before computing similarities. Memorization is quantified at the level of individual generated samples. For each generated sample $\tilde{x}$, we compute its two nearest training neighbors and define distances d_1 and d_2. A sample is considered memorized if $\frac{d_1}{d_2} < \tau$ for a range of thresholds $\tau \in \{1/6, 1/4, 1/3, 1/2, 2/3\}$. This yields both per-sample and per-training-point memorization statistics. Local data density is estimated using nearest-neighbor statistics computed solely on the training set. For each training sample, we compute the distances to the k nearest neighbors for $k \in \{2, 5, 10, 20, 50\}$ and average over these distances as defined in Equation (18). These statistics serve as proxies for local sparsity and are later correlated with memorization behavior.

C.1. CIFAR-10 Experiments

We construct the CIFAR-10 dataset by combining the original training and test splits (60,000 images total), and then randomly partitioning them into three equal subsets (train/validation/test), each containing approximately 20,000 images. All images are represented as 32×32 RGB tensors normalized to $[0, 1]$. No additional preprocessing or augmentation is applied.

C.2. MNIST + CIFAR-10 Experiments

To study memorization under controlled differences in local data density, we construct a combined dataset from CIFAR-10 and MNIST, which exhibit markedly different intrinsic complexities. CIFAR-10 images are highly variable, while MNIST digits occupy a much more concentrated region of image space. This setup allows us to induce systematic differences in local sparsity across classes.

We merge the original training and test splits of both datasets. To ensure compatibility, MNIST images are resized to 32×32 and converted to three channels by duplicating the grayscale channel, resulting in RGB tensors consistent with CIFAR-10. CIFAR-10 images are used without modification. All images are represented as 32×32 RGB tensors normalized to $[0, 1]$, and no additional preprocessing or augmentation is applied.

To isolate dataset-level effects, we ignore the original class labels and instead assign a binary label indicating the dataset of origin (CIFAR-10 or MNIST). The two datasets are then balanced by subsampling the larger dataset such that both contribute an equal number of samples. The resulting dataset therefore consists of an equal mixture of CIFAR-10 and MNIST images.

C.3. CelebA Experiments

This dataset consists of high-resolution face images together with 40 binary attributes per image. All images are resized to 32×32 pixels, and converted to grayscale. The resulting images are represented as single-channel tensors normalized to $[0, 1]$. No additional data augmentation is applied.