

Diffusion Model: A Detailed Tutorial

May, 2025

Contents

1	SDE Representation of the Forward Process in DDPM	2
2	SDE Representation of the Reverse Process in DDPM	2
3	Incorporating a Neural Network	3
4	Discretizing the Reverse Process: A Stepwise View	4
5	Interpreting Diffusion Models via the Score Function	5
6	The DPS Algorithm: Posterior-Consistent Generation	5
6.1	Algorithmic Rationale	5
6.2	Practical Implementation	6
7	DDIM: A Higher-Level Perspective on DDPM	7
8	Concrete Implementation of DDPM	7
8.1	Training Procedure	7
8.2	Inference (Sampling) Procedure	7
8.3	Extension to Text-to-Image Generation	8
9	DreamFusion: From 2-D to 3-D Generation	8
9.1	Conceptual Overview	8
9.2	Density Distillation in Parameter Space	8
9.3	Practical Pipeline (Stable DreamFusion)	9
10	ProlificDreamer: Optimizing Parameter <i>Distributions</i>	9
10.1	Principle Overview	9
10.1.1	Variational Score Distillation (VSD)	9
10.1.2	Updating μ via Particle-Based Variational Inference	9
10.2	Practical Workflow (Pseudo-Code)	10

1 SDE Representation of the Forward Process in DDPM

During the noise-adding stage of a **DDPM**, every time-step applies the following discrete Markov chain:

$$x_i = \sqrt{1 - \beta_i} x_{i-1} + \sqrt{\beta_i} \varepsilon_{i-1}, \quad i = 1, \dots, N. \quad (1)$$

To turn this into a continuous process, we let the discrete interval Δt shrink to zero-equivalently, we consider the limit $N \rightarrow \infty$ of the Markov chain.

Before taking the limit, introduce an auxiliary noise scale $\{\tilde{\beta}_i = N \beta_i\}_{i=1}^N$ and rewrite

$$x_i = \sqrt{1 - \frac{\tilde{\beta}_i}{N}} x_{i-1} + \sqrt{\frac{\tilde{\beta}_i}{N}} \varepsilon_{i-1}, \quad i = 1, \dots, N. \quad (2)$$

As $N \rightarrow \infty$, the sequence $\{\tilde{\beta}_i\}_{i=1}^N$ becomes a continuous schedule $\beta(t)$ on $t \in [0, 1]$. Set

$$\Delta t = \frac{1}{N}.$$

At each Δt , the continuous functions $\beta(t)$, $x(t)$, $\varepsilon(t)$ coincide with their discrete counterparts:

$$\beta\left(\frac{i}{N}\right) = \tilde{\beta}_i, \quad x\left(\frac{i}{N}\right) = x_i, \quad \varepsilon\left(\frac{i}{N}\right) = \varepsilon_i.$$

For $t \in \{0, \frac{1}{N}, \dots, \frac{N-1}{N}\}$ and $t + \Delta t = t + \frac{1}{N}$, we rewrite the update with continuous notation:

$$x(t + \Delta t) = \sqrt{1 - \beta(t + \Delta t) \Delta t} x(t) + \sqrt{\beta(t + \Delta t) \Delta t} \varepsilon(t) \quad (3)$$

$$\approx x(t) - \frac{1}{2} \beta(t + \Delta t) \Delta t x(t) + \sqrt{\beta(t + \Delta t) \Delta t} \varepsilon(t) \quad (4)$$

$$\approx x(t) - \frac{1}{2} \beta(t) \Delta t x(t) + \sqrt{\beta(t) \Delta t} \varepsilon(t). \quad (5)$$

(The second line holds when $\Delta t \ll 1$.) Thus

$$x(t + \Delta t) - x(t) \approx -\frac{1}{2} \beta(t) \Delta t x(t) + \sqrt{\beta(t) \Delta t} \varepsilon(t).$$

Letting $\Delta t \rightarrow 0$ gives the stochastic differential equation (SDE)

$$\boxed{dx = -\frac{1}{2} \beta(t) x dt + \sqrt{\beta(t)} dw}. \quad (6)$$

2 SDE Representation of the Reverse Process in DDPM

Using probabilistic language, Eq. (6) can be generalized to the stochastic differential equation (SDE)

$$dx = f_t(x) dt + g_t dw, \quad (7)$$

which likewise describes the forward noise-adding Markov chain. Whether the chain is *linear* depends on whether $f_t(x)$ is linear in x . The corresponding one-step conditional density is

$$p(x_{t+\Delta t} | x_t) = \mathcal{N}\left(x_{t+\Delta t}; x_t + f_t(x_t) \Delta t, g_t^2 \Delta t I\right) \propto \exp\left(-\frac{\|x_{t+\Delta t} - x_t - f_t(x_t) \Delta t\|^2}{2 g_t^2 \Delta t}\right). \quad (8)$$

(The constant normalization factor is omitted.)

Following the DDPM philosophy—“learning to build by watching demolition”—we ultimately wish to find the reverse density $p(x_t | x_{t+\Delta t})$. Applying Bayes’ rule,

$$\begin{aligned} p(x_t | x_{t+\Delta t}) &= \frac{p(x_{t+\Delta t} | x_t) p(x_t)}{p(x_{t+\Delta t})} \\ &\propto \exp\left(-\frac{\|x_{t+\Delta t} - x_t - f_t(x_t) \Delta t\|^2}{2 g_t^2 \Delta t} + \log p(x_t) - \log p(x_{t+\Delta t})\right). \end{aligned} \quad (9)$$

Because Δt is small, $p(x_{t+\Delta t} | x_t)$ is appreciable only when $x_{t+\Delta t}$ is close to x_t ; the same is true for the reverse density. Hence we expand

$$\log p(x_{t+\Delta t}) \approx \log p(x_t) + (x_{t+\Delta t} - x_t)^\top \nabla_{x_t} \log p(x_t) + \Delta t \partial_t \log p(x_t).$$

Substituting and collecting terms,

$$p(x_t | x_{t+\Delta t}) \propto \exp\left(-\frac{\|x_{t+\Delta t} - x_t - [f_t(x_t) - g_t^2 \nabla_{x_t} \log p(x_t)] \Delta t\|^2}{2 g_t^2 \Delta t} + \mathcal{O}(\Delta t)\right). \quad (10)$$

Letting $\Delta t \rightarrow 0$ gives

$$p(x_t | x_{t+\Delta t}) \propto \exp\left(-\frac{\|x_{t+\Delta t} - x_t - [f_{t+\Delta t}(x_{t+\Delta t}) - g_{t+\Delta t}^2 \nabla_{x_{t+\Delta t}} \log p(x_{t+\Delta t})] \Delta t\|^2}{2 g_{t+\Delta t}^2 \Delta t}\right). \quad (11)$$

Hence $p(x_t | x_{t+\Delta t})$ is approximately Gaussian with mean

$$x_{t+\Delta t} - \left[f_{t+\Delta t}(x_{t+\Delta t}) - g_{t+\Delta t}^2 \nabla_{x_{t+\Delta t}} \log p(x_{t+\Delta t}) \right] \Delta t,$$

and covariance $g_{t+\Delta t}^2 \Delta t I$. Taking $\Delta t \rightarrow 0$ recovers the reverse-time SDE

$$dx = [f_t(x) - g_t^2 \nabla_x \log p(x)] dt + g_t dw. \quad (12)$$

For the linear forward Markov chain in Eq. (1), $f_t(x) = -\frac{1}{2}\beta(t)x$ and $g_t = \sqrt{\beta(t)}$, yielding

$$\boxed{dx = \left[-\frac{1}{2}\beta(t)x - \beta(t)\nabla_x \log p(x)\right] dt + \sqrt{\beta(t)} dw}. \quad (13)$$

3 Incorporating a Neural Network

Given the reverse-time SDE in Eq. (13), the noise schedule $\beta(t)$ is known at every step; to complete the *generation* (or “building-up”) process we need only the score $\nabla_x \log p(x)$. Here $p(x)$ denotes the marginal distribution of x at time t . For a *linear* forward Markov process this marginal can be written in closed form, but doing so requires an average over *all* training samples x_0 , which is computationally expensive and does not generalize well. We therefore train a neural network to *directly* approximate $\nabla_x \log p(x)$.

The diffusion process specifies the forward transition density $p(x_{t+\Delta t} | x_t)$. Integrating these infinitesimal transitions in sequence gives

$$p(x_t | x_0) = \lim_{\Delta t \rightarrow 0} \int \cdots \int p(x_t | x_{t-\Delta t}) p(x_{t-\Delta t} | x_{t-2\Delta t}) \cdots p(x_{\Delta t} | x_0) dx_{t-\Delta t} \cdots dx_{\Delta t}. \quad (14)$$

When the forward chain is linear, the expression above admits a closed-form solution (not guaranteed if the chain is nonlinear). Consequently the marginal at time t is

$$p(x_t) = \int p(x_t | x_0) p(x_0) dx_0 = \mathbb{E}_{x_0}[p(x_t | x_0)], \quad (15)$$

and

$$\nabla_{x_t} \log p(x_t) = \frac{\mathbb{E}_{x_0}[p(x_t | x_0) \nabla_{x_t} \log p(x_t | x_0)]}{\mathbb{E}_{x_0}[p(x_t | x_0)]}. \quad (16)$$

Lemma 1 (Weighted MSE minimizer). *Let $x \in \mathbb{R}^d$ be a random vector with $\mathbb{E}\|x\|^2 < \infty$, and let $w \geq 0$ be a non-negative random weight on the same space with $\mathbb{E}[w] > 0$. For any fixed vector $y \in \mathbb{R}^d$, define the weighted mean-squared error*

$$J(y) = \mathbb{E}[w \|y - x\|^2].$$

Then $J(y)$ is strictly convex in y and attains its unique minimizer at

$$y^* = \frac{\mathbb{E}[w x]}{\mathbb{E}[w]}.$$

When $w \equiv 1$, this reduces to the classical result $y^ = \mathbb{E}[x]$.*

Choose

$$w = p(x_t | x_0), \quad x = \nabla_{x_t} \log p(x_t | x_0), \quad y = s_\theta(x_t, t),$$

and define the loss

$$\mathcal{L}(s_\theta) = \mathbb{E}_{x_0} [p(x_t | x_0) \|s_\theta(x_t, t) - \nabla_{x_t} \log p(x_t | x_0)\|^2]. \quad (17)$$

By the lemma, this loss is minimized at

$$s_\theta^*(x_t, t) = \frac{\mathbb{E}_{x_0} [p(x_t | x_0) \nabla_{x_t} \log p(x_t | x_0)]}{\mathbb{E}_{x_0} [p(x_t | x_0)]} = \nabla_{x_t} \log p(x_t). \quad (18)$$

Since the denominator merely rescales the loss, we drop it for simplicity. Expanding the outer expectation over x_0 yields

$$\mathcal{L}(s_\theta) = \int \mathbb{E}_{x_0|x_t} [\|s_\theta(x_t, t) - \nabla_{x_t} \log p(x_t | x_0)\|^2] p(x_t) dx_t = \mathbb{E}_{x_0, x_t \sim p(x_t|x_0) \tilde{p}(x_0)} [\|s_\theta(x_t, t) - \nabla_{x_t} \log p(x_t | x_0)\|^2] \quad (19)$$

Forward-process reparameterization. From the DDPM forward dynamics one can write x_t in closed form:

$$x_t = \alpha_t x_{t-1} + \beta_t \varepsilon_t = \dots = (\alpha_t \dots \alpha_1) x_0 + \sum_{k=1}^t (\alpha_t \dots \alpha_{k+1}) \beta_k \varepsilon_k.$$

Because $(\alpha_t \dots \alpha_1)^2 + (\alpha_t \dots \alpha_2)^2 \beta_1^2 + \dots + \beta_t^2 = 1$, we may write

$$x_t = \tilde{\alpha}_t x_0 + \tilde{\beta}_t \varepsilon, \quad \tilde{\alpha}_t^2 + \tilde{\beta}_t^2 = 1,$$

with $p(x_t | x_0) = \mathcal{N}(x_t; \tilde{\alpha}_t x_0, \tilde{\beta}_t^2 I)$.

Hence

$$\nabla_{x_t} \log p(x_t | x_0) = -\frac{1}{\tilde{\beta}_t^2} (x_t - \tilde{\alpha}_t x_0) = -\frac{1}{\tilde{\beta}_t} \varepsilon.$$

Let $s_\theta(x_t, t) = -\varepsilon_\theta(x_t, t)/\tilde{\beta}_t$; dropping the constant factor $1/\tilde{\beta}_t^2$ gives the final training loss

$$\boxed{\mathcal{L}(\varepsilon_\theta) = \mathbb{E}_{x_0 \sim \tilde{p}(x_0), \varepsilon \sim \mathcal{N}(0, I)} [\|\varepsilon_\theta(\tilde{\alpha}_t x_0 + \tilde{\beta}_t \varepsilon, t) - \varepsilon\|^2]}. \quad (20)$$

Minimizing this loss forces the network ε_θ to approximate the true score $\nabla_{x_t} \log p(x_t)$ at *every* time step t .

4 Discretizing the Reverse Process: A Stepwise View

Sections 1 and 2 showed that a trained DDPM can be interpreted as a stochastic process that models the *reverse* (denoising) dynamics. Here we discretize those continuous dynamics and make explicit how to sample x_{t-1} from a given x_t .

By Bayes' theorem

$$p(x_{t-1} | x_t) = \frac{p(x_t | x_{t-1}) p(x_{t-1})}{p(x_t)}. \quad (21)$$

Because the marginals $p(x_{t-1})$ and $p(x_t)$ are unknown in closed form, this expression is not directly usable (the continuous-time derivation in Section 2 sidestepped them by taking $\Delta t \rightarrow 0$).

Instead, condition on the (unknown) clean image x_0 :

$$p(x_{t-1} | x_t, x_0) = \frac{p(x_t | x_{t-1}, x_0) p(x_{t-1} | x_0)}{p(x_t | x_0)}. \quad (22)$$

Now every factor is known analytically, yielding

$$p(x_{t-1} | x_t, x_0) = \mathcal{N}\left(x_{t-1}; \frac{\alpha_t \tilde{\beta}_{t-1}^2}{\tilde{\beta}_t^2} x_t + \frac{\bar{\alpha}_{t-1} \beta_t^2}{\tilde{\beta}_t^2} x_0, \frac{\tilde{\beta}_{t-1}^2}{\tilde{\beta}_t^2} I\right). \quad (23)$$

Because x_0 is unavailable during inference, we replace it by a learned predictor $\hat{\mu}(x_t)$. If $\hat{\mu}(x_t)$ is trained with the loss $\|x_0 - \hat{\mu}(x_t)\|^2$, then

$$p(x_{t-1} | x_t) \approx p(x_{t-1} | x_t, x_0 = \hat{\mu}(x_t)) = \mathcal{N}\left(x_{t-1}; \frac{\alpha_t \bar{\beta}_{t-1}^2}{\bar{\beta}_t^2} x_t + \frac{\bar{\alpha}_{t-1} \beta_t^2}{\bar{\beta}_t^2} \hat{\mu}(x_t), \frac{\bar{\beta}_{t-1}^2}{\bar{\beta}_t^2} I\right).$$

Because $x_t = \bar{\alpha}_t x_0 + \bar{\beta}_t \varepsilon$, one may solve for x_0 as $x_0 = \frac{1}{\bar{\alpha}_t}(x_t - \bar{\beta}_t \varepsilon)$. This motivates

$$\hat{\mu}(x_t) = \frac{1}{\bar{\alpha}_t} \left(x_t - \bar{\beta}_t \varepsilon_\theta(x_t, t) \right),$$

where ε_θ is the usual noise-predicting UNet.

Training ε_θ with

$$\|x_0 - \hat{\mu}(x_t)\|^2 = \frac{\bar{\beta}_t^2}{\bar{\alpha}_t^2} \left\| \varepsilon - \varepsilon_\theta(\bar{\alpha}_t x_0 + \bar{\beta}_t \varepsilon, t) \right\|^2$$

reduces (up to a scalar) to the simplified DDPM loss.

Substituting $\hat{\mu}(x_t)$ back yields

$$p(x_{t-1} | x_t) \approx \mathcal{N}\left(x_{t-1}; \frac{1}{\alpha_t} \left(x_t - \frac{\beta_t^2}{\bar{\beta}_t} \varepsilon_\theta(x_t, t) \right), \frac{\bar{\beta}_{t-1}^2 \beta_t^2}{\bar{\beta}_t^2} I\right),$$

so that a single sampling step is

$$x_{t-1} = \frac{1}{\alpha_t} \left(x_t - \frac{\beta_t^2}{\bar{\beta}_t} \varepsilon_\theta(x_t, t) \right) + \frac{\bar{\beta}_{t-1} \beta_t}{\bar{\beta}_t} \varepsilon, \quad \varepsilon \sim \mathcal{N}(0, I). \quad (24)$$

5 Interpreting Diffusion Models via the Score Function

The *score function* of a density is defined as its log-likelihood gradient, $s(x) = \nabla_x \log p(x)$, pointing in the direction of steepest increase in probability. In Langevin dynamics and diffusion models, this score acts as the drift term that “pulls” noisy samples back toward high-density regions, as schematically illustrated below.

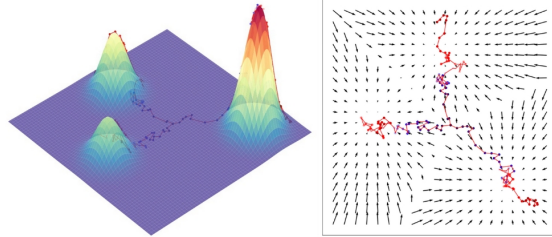


Figure 1: Intuition: score-based drift guides noisy particles toward the data manifold.

Accordingly, the reverse process trains the network to approximate the score of each noisy image x_t , effectively teaching it how to guide particles from $p(x_{t+1})$ back to $p(x_t)$.

6 The DPS Algorithm: Posterior-Consistent Generation

6.1 Algorithmic Rationale

Armed with the preceding background, we now examine the **DPS (Denoising Posterior Sampling)** algorithm in a practical setting.

In an *unconditional* diffusion model the only guidance is the prior score $\nabla_x \log p(x)$; the learned “force field” merely pulls noisy samples toward the natural-image manifold. In real applications—say, image denoising—we need the final output to be both *realistic* (prior fidelity) and *consistent* with a

noisy observation y (data fidelity). Hence we upgrade the marginal $p(x)$ at every step to a conditional $p(x | y)$. In the reverse SDE, the network must now predict $\nabla_x \log p(x | y)$ instead of $\nabla_x \log p(x)$.

Bayes' rule gives

$$\nabla_x \log p(x | y) = \nabla_x \log p(x) + \nabla_x \log p(y | x).$$

The prior term can be pretrained as usual, but $\nabla_x \log p(y | x)$ lacks a closed form because the likelihood is time-dependent and we only know the relation between y and the clean image x_0 .

Assume the observation model $y = A(x_0) + n$, with A an imaging operator and n additive noise. Because the likelihood $p(y | x_t)$ has no analytic form, we relate it to the known $p(y | x_0)$ via conditional independence:

$$p(y | x_t) = \int p(y | x_0, x_t) p(x_0 | x_t) dx_0 = \int p(y | x_0) p(x_0 | x_t) dx_0 = \mathbb{E}_{x_0 \sim p(x_0 | x_t)}[p(y | x_0)].$$

Applying Jensen's inequality,

$$p(y | x_t) \approx p(y | \hat{x}_0), \quad \hat{x}_0 := \mathbb{E}[x_0 | x_t].$$

Because the forward process satisfies $x_t = \sqrt{\bar{\alpha}} x_0 + \sqrt{1 - \bar{\alpha}} z$, one finds

$$\hat{x}_0(x_t) = \frac{1}{\sqrt{\bar{\alpha}}} \left(x_t + (1 - \bar{\alpha}) \nabla_{x_t} \log p_t(x_t) \right).$$

Assuming Gaussian observation noise

$$p(y | x_0) \propto \exp\left(-\frac{1}{2\sigma^2} \|y - A(x_0)\|_2^2\right),$$

we obtain

$$\nabla_{x_t} \log p_t(y | x_t) \approx -\frac{1}{\sigma^2} \nabla_{x_t} \|y - A(\hat{x}_0(x_t))\|_2^2.$$

Hence the *posterior* score is approximated by

$$\nabla_{x_t} \log p_t(x_t | y) \approx s_\theta^*(x_t, t) - \rho \nabla_{x_t} \|y - A(\hat{x}_0)\|_2^2, \quad \rho := \frac{1}{\sigma^2}.$$

6.2 Practical Implementation

1. **Initialization.** Start the reverse chain with pure noise $x_N \sim \mathcal{N}(0, I)$.
2. **Backward iteration.** For $i = N-1, \dots, 0$:
 - (i) Compute the prior score with the pretrained model $\hat{s} = s_\theta(x_i, i) \approx \nabla_{x_i} \log p_t(x_i)$.
 - (ii) **Tweedie update** (estimate a clean image)

$$\hat{x}_0 = \frac{1}{\sqrt{\bar{\alpha}_i}} (x_i + (1 - \bar{\alpha}_i) \hat{s}).$$

- (iii) **Prior-only DDPM step**

$$x'_{i-1} = \sqrt{\frac{\bar{\alpha}_{i-1}(1 - \alpha_i)}{1 - \bar{\alpha}_i}} x_i + \sqrt{\frac{\bar{\alpha}_{i-1}\beta_i}{1 - \bar{\alpha}_i}} \hat{x}_0 + \tilde{\sigma}_i z, \quad z \sim \mathcal{N}(0, I).$$

- (iv) **Likelihood correction** (Gaussian case)

$$x_{i-1} = x'_{i-1} - \zeta_i A^\top (A(\hat{x}_0) - y),$$

where ζ_i trades off prior versus data fidelity.

3. **Output.** After the loop, return the reconstruction \hat{x}_0 —now consistent with both the prior (realism) and the observation y (data fidelity).

7 DDIM: A Higher-Level Perspective on DDPM

Why model diffusion with the *solution* of an SDE—i.e. a *stochastic* process? Because an SDE captures the key feature of DDPM denoising: given the previous state, the next state is *not* deterministic; only its *distribution* is specified. Suppose all images possessing a certain attribute follow a latent distribution q . We want a model that, when fed a noisy image x_t at noise level t , predicts *the distribution of x_{t-1}* , not a single value. By sampling from that distribution at each step we eventually obtain a random draw from q . If a large noise level T converts any image to roughly standard normal, then by T steps of prediction and sampling we can transform pure Gaussian noise into a valid sample from q .

Once the forward relation $x_t \leftrightarrow x_0$ is available in closed form, the step-by-step noising is unnecessary; the time parameter t merely controls the noise intensity. This observation leads to DDIM: because the *result* does not depend on $p(x_t | x_{t-1})$, we can drop the “build-and-demolish” construction entirely.

In principle, even without an explicit $p(x_t | x_{t-1})$, the conditional $p(x_{t-1} | x_t, x_0)$ is solvable; indeed, the solution set is larger and easier to characterize. All that is required is the *marginal-consistency* condition

$$\int p(x_{t-1} | x_t, x_0) p(x_t | x_0) dx_t = p(x_{t-1} | x_0).$$

With undetermined coefficients we can solve this directly. More generally, assume

$$p(x_{t-1} | x_t, x_0) = \mathcal{N}(x_{t-1}; \kappa_t x_t + \lambda_t x_0, \sigma_t^2 I),$$

with $\kappa_t, \lambda_t, \sigma_t$ to be determined. Using $p(x_{t-1} | x_0)$ and $p(x_t | x_0)$ one obtains a *family* of solutions parameterized by the free variance σ_t . Training is unaffected (the saved model is unchanged), but generation now has a tunable parameter σ_t —the key novelty introduced by DDIM.

In the building-demolishing metaphor, we know what the fully demolished building looks like ($p(x_t | x_0)$ and $p(x_{t-1} | x_0)$), but not how each individual plank is removed ($p(x_t | x_{t-1})$). If x_t lets us estimate x_0 , then teaching the model, given (x_0, x_t) , to recover the intermediate state *is* to learn every reverse step $p(x_{t-1} | x_t)$ —in other words, to “rebuild” the house one floor at a time without ever specifying the original demolition plan.

8 Concrete Implementation of DDPM

8.1 Training Procedure

- Sample an image x_0 from the training set ($x_0 \sim q(x_0)$); Draw a random time step $t \sim \text{Uniform}(1, \dots, T)$;
- Sample Gaussian noise $\varepsilon \sim \mathcal{N}(0, \mathbf{I})$;
- Compute the loss

$$\text{loss} = \left\| \varepsilon - \varepsilon_\theta(\sqrt{\bar{\alpha}_t} x_0 + \sqrt{1 - \bar{\alpha}_t} \varepsilon, t) \right\|^2,$$

where ε_θ is the UNet-based denoising network;

- Back-propagate the loss and update the model parameters; repeat until convergence.

8.2 Inference (Sampling) Procedure

- Draw an initial latent $x_T \sim \mathcal{N}(0, \mathbf{I})$;
- For $t = T, T - 1, \dots, 1$:
 - Sample fresh noise $\varepsilon \sim \mathcal{N}(0, \mathbf{I})$;
 - Use Eq. (24) to obtain x_{t-1} from x_t .
- The final output is x'_0 .

8.3 Extension to Text-to-Image Generation

To turn a DDPM into a text-to-image diffusion model, replace each self-attention block in the UNet with a *cross-attention* block, feeding the encoded text prompt y as \mathbf{K} and \mathbf{V} while retaining the latent features as \mathbf{Q} . The training loss becomes

$$\text{loss} = \left\| \varepsilon - \varepsilon_\theta(\sqrt{\bar{\alpha}_t} x_0 + \sqrt{1 - \bar{\alpha}_t} \varepsilon, t, y) \right\|^2.$$

9 DreamFusion: From 2-D to 3-D Generation

Diffusion models have achieved impressive text-to-*image* generation, yet text-to-*3-D* remains difficult because large text-shape pairs are scarce. *DreamFusion: Text-to-3D Using 2-D Diffusion* remedies this by using a *pre-trained text-to-image* diffusion model to supervise a 3-D generator.

9.1 Conceptual Overview

Let θ denote the parameters of a volume renderer $g(\cdot)$ (e.g. a Gaussian NeRF). Render $x = g(\theta)$, feed it into a frozen U-Net, and define the diffusion loss

$$\mathcal{L}_{\text{Diff}}(\varphi, x = g(\theta)) = \mathbb{E}_{t, \varepsilon} [\| \hat{\varepsilon}_\varphi(\alpha_t g(\theta) + \sigma_t \varepsilon; y, t) - \varepsilon \|^2],$$

where $\hat{\varepsilon}_\varphi$ is the frozen U-Net predictor. If $g(\theta)$ produces photorealistic pixels, the U-Net predicts noise well and the loss is low; otherwise it is high.

With $z_t = \alpha_t x + \sigma_t \varepsilon$ and $x = g(\theta)$,

$$\nabla_\theta \mathcal{L}_{\text{Diff}} = \mathbb{E}_{t, \varepsilon} [2(\hat{\varepsilon}_\varphi(z_t) - \varepsilon) \partial_{z_t} \hat{\varepsilon}_\varphi(z_t) \partial_x z_t \partial_\theta x].$$

Here $\partial_{z_t} \hat{\varepsilon}_\varphi$ (the U-Net Jacobian) is enormous; propagating it would require a full backward pass *per* pixel sample and yields unstable gradients at high noise levels.

Therefore DreamFusion *drops* the Jacobian, retaining only the residual $\hat{\varepsilon}_\varphi(z_t) - \varepsilon$. The gradient becomes

$$\nabla_\theta \mathcal{L}_{\text{SDS}} = \mathbb{E}_{t, \varepsilon} [w(t) (\hat{\varepsilon}_\varphi(z_t; y, t) - \varepsilon) \partial_\theta x],$$

the celebrated **score-distillation sampling (SDS)** gradient.

Crucially, SDS is **not** a heuristic; it performs *probability-density distillation* in parameter space: updating θ with the SDS gradient equals minimizing a *weighted KL divergence* between (1) the noisy distribution of the rendered image and (2) the true diffusion distribution at the same noise level t .

9.2 Density Distillation in Parameter Space

KL divergence. For densities $p(z)$ and $q(z)$,

$$\text{KL}(p\|q) = \int p(z) \log \frac{p(z)}{q(z)} dz = \mathbb{E}_{z \sim p} [\log p(z) - \log q(z)].$$

If q encodes p , the extra code length over the optimal $H(p)$ is exactly $\text{KL}(p\|q)$.

Equivalence to SDS. Let $q(z_t | x) = \mathcal{N}(z_t; \alpha_t x, \sigma_t^2 I)$ with $x = g(\theta)$, and $p_\phi(z_t | y)$ be the true diffusion distribution at level t . One finds

$$\nabla_\theta \text{KL}(q\|p_\phi) = \mathbb{E}_{t, \varepsilon} [w(t) (\epsilon_\phi(z_t) - \varepsilon) \partial_\theta x],$$

identical to $\nabla_\theta \mathcal{L}_{\text{SDS}}$.

9.3 Practical Pipeline (Stable DreamFusion)

Each step:

1. **Rendering.** Render a low-resolution image $x = g(\theta)$ ($[B, 3, H, W]$), then upsample.
2. **Forward noise.** Encode the high-res image with the VAE of Stable Diffusion; sample $t \sim \text{Uniform}(1, \dots, T)$ and add noise.
3. **Denoising.** Run the noisy latent through the frozen U-Net to obtain $\hat{\epsilon}$.
4. **Back-prop SDS.** Because the U-Net is frozen, stop its gradient and compute

```

1 w      = (1 - self.alphas[t])                # weight w(t)
2 grad    = w[:, None, None, None] * (noise_pred - noise)
3 target  = (latents - grad).detach()           # stop gradient
4 loss_sds = 0.5 * F.mse_loss(
5         latents, target, reduction="sum") / batch_size

```

5. **Update θ .** Back-propagate `loss_sds` through the NeRF MLP only; iterate until convergence.

10 ProlificDreamer: Optimizing Parameter *Distributions*

The previous section (DreamFusion) trains a *single* NeRF parameter set θ via diffusion and the SDS loss, but such a point estimate can suffer from *low diversity*. Inspired by 2-D diffusion, where one first learns a *distribution* over images and then samples from it, we ask: can we likewise learn a *distribution* μ over 3-D parameters θ and then sample a diverse θ from μ ? This is the core idea behind **ProlificDreamer**.

10.1 Principle Overview

10.1.1 Variational Score Distillation (VSD)

For a text prompt y there exists a distribution of all 3-D scenes consistent with that prompt. Denote the distribution of NeRF parameters by $\mu(\theta | y)$. Rendering with viewpoint c yields $x_0 = g(\theta, c)$. Let $q_0^\mu(x_0 | c, y)$ be the image distribution obtained by sampling $\theta \sim \mu(\cdot | y)$ and rendering at c . Let $p_0(x_0 | y)$ be the distribution produced by a frozen text-to-image diffusion model conditioned on y .

We wish to minimize their KL divergence:

$$\min_{\mu} D_{\text{KL}}(q_0^\mu(x_0 | y) \| p_0(x_0 | y)),$$

a *classical variational-inference* objective.

To ensure closeness *at every noise level*, we instead minimize the weighted KL along the entire noising trajectory:

$$\mu^* = \arg \min_{\mu} \mathbb{E}_{t,c} \left[\frac{\sigma_t}{\alpha_t} w(t) D_{\text{KL}}(q_t^\mu(x_t | c, y) \| p_t(x_t | y)) \right]. \quad (25)$$

10.1.2 Updating μ via Particle-Based Variational Inference

Treat n parameter sets $\{\theta_i\}_{i=1}^n$ as *particles* representing μ ($n=4$ in the original paper). Gradient descent with step size $\eta \rightarrow 0$ leads to an ODE:

$$\dot{\theta}_\tau = -\nabla_{\theta} L(\theta_\tau),$$

known here as the *Wasserstein gradient flow of VSD*. Starting from $\theta_0 \sim \mu_0(\theta | y)$ and integrating the ODE yields the optimal distribution μ_τ as $\tau \rightarrow \infty$.

Define

$$\frac{d\theta_\tau}{d\tau} = -\mathbb{E}_{t,\varepsilon,c} \left[\omega(t) \left(-\sigma_t \nabla_{\mathbf{x}_t} \log p_t(\mathbf{x}_t | y) + \sigma_t \nabla_{\mathbf{x}_t} \log q_t^{\mu_\tau}(\mathbf{x}_t | c, y) \right) \frac{\partial g(\theta_\tau, c)}{\partial \theta_\tau} \right].$$

Real-image scores use the frozen diffusion model $\epsilon_{\text{pretrain}}$; rendered-image scores use a learnable network ϵ_ϕ (a lightweight UNet or LoRA-tuned clone of the pretrained network). Alternating updates yield

$$\nabla_{\theta} \mathcal{L}_{\text{VSD}}(\theta) = \mathbb{E}_{t, \varepsilon, c} \left[\omega(t) (\epsilon_{\text{pretrain}}(\mathbf{x}_t, t, y) - \epsilon_\phi(\mathbf{x}_t, t, c, y)) \frac{\partial g(\theta, c)}{\partial \theta} \right].$$

SDS is the special case where μ is a Dirac delta.

10.2 Practical Workflow (Pseudo-Code)

Inputs: number of particles n , prompt y , frozen score $\epsilon_{\text{pretrain}}$, learning rates η_1, η_2 , renderer $g(\theta, c)$, noise schedule $\{\alpha_t, \sigma_t, \omega(t)\}$.

Initialization: $\{\theta_0^{(i)}\}_{i=1}^n \sim \text{InitPrior}$, $\phi_0 \leftarrow \text{InitWeights}$.

Main Loop: for $k = 0, \dots$ until convergence

1. Sample particle index and camera: $i_k \sim \text{Uniform}\{1, \dots, n\}$, $c_k \sim p(c)$.
2. Render clean image: $x_0^{(k)} = g(\theta_k^{(i_k)}, c_k)$.
3. Noise it: $t_k \sim \text{U}(0, 1)$, $\varepsilon_k \sim \mathcal{N}(0, I)$, $x_{t_k}^{(k)} = \alpha_{t_k} x_0^{(k)} + \sigma_{t_k} \varepsilon_k$.
4. Compute real and rendered scores: $\hat{\varepsilon}_k^{\text{real}} = \epsilon_{\text{pretrain}}(x_{t_k}^{(k)}, t_k, y)$, $\hat{\varepsilon}_k^{\text{rend}} = \epsilon_{\phi_k}(x_{t_k}^{(k)}, t_k, c_k, y)$.
5. Update geometry:

$$\theta_{k+1}^{(i_k)} = \theta_k^{(i_k)} - \eta_1 \omega(t_k) (\hat{\varepsilon}_k^{\text{real}} - \hat{\varepsilon}_k^{\text{rend}}) \frac{\partial g(\theta_k^{(i_k)}, c_k)}{\partial \theta^{(i_k)}}.$$

6. Update ϕ : minimize $\|\epsilon_{\phi_k}(x_{t_k}^{(k)}, t_k, c_k, y) - \varepsilon_k\|_2^2$ with step η_2 .

Return the final particle set $\{\theta_{\text{final}}^{(i)}\}_{i=1}^n$ and score network ϕ_{final} .