

3D Reconstruction with Fast Dipole Sums

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Abstract

Reconstructing 3D scenes from multi-view images has always been a challenging problem in computer vision and computer graphics. Traditional methods like structure from motion and multi-view stereo have been widely used for pose estimation and dense point cloud reconstruction. However, these methods have limited ability to reconstruct complex scenes with fine details. Recently, since the introduction of neural radiance fields (NeRF), volumetric neural rendering has shown great promise in reconstructing complex scenes with high fidelity. To accurately reconstruct scene geometry, other works have also proposed ways to directly model the signed-distance function or occupancy of a scene. However, these methods are often slow to train and cannot effectively leverage known scene information.

In this thesis, we propose a novel point-based representation that combines the efficiency of point clouds with the expressiveness of neural rendering. Point clouds are particularly appealing as a scene representation for rendering tasks, as they are the natural output of many 3D sensing modalities, including structure from motion, multi-view stereo, and lidar. They also come with a rich library of geometric queries. In our work, we utilize point clouds to efficiently reconstruct 3D scenes by using the generalized winding number as a proxy for the scene occupancy and by interpolating per-point neural features with appropriate kernels. We leverage the Barnes-Hut approximation and fast dipole sums to perform fast winding number queries and feature interpolation, as well as logarithmic complexity backpropagation for efficient differentiable rendering. We empirically show that our method consistently outperforms existing methods in both reconstruction quality and efficiency on a wide range of real-world scenes.

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Chapter 1

Related work

Things to cite here: - Nick's paper, Belkin paper, on point cloud geometry - SIGGRAPH 2023 paper on fixing normal orientations - Point-NeRF - CMU paper on using point clouds - GeoNeuS - Pointersect - Gaussian splatting - Kaleidoscopic, other stereo that produce area weights - Alec's winding number papers - PSR and SPSR - Neural surface papers - Differentiable PSR

Chapter 2

Volumetric neural rendering

2.1 Volumetric light transport background

We first introduce the basic concepts of volumetric light transport using principles from classical volume rendering [9], which will serve as the foundation for our rendering framework.

2.1.1 Radiative transfer equation

The *radiative transfer equation* (RTE) describes the propagation of light in a medium. It is a partial differential equation that models the change in radiance along a ray as it travels through the medium. Assuming an emissive medium with no scattering, the RTE can be written as:

$$dL(\mathbf{x}, \vec{\omega}) = -\sigma(\mathbf{x}, \vec{\omega})L(\mathbf{x}, \vec{\omega})dz + \sigma(\mathbf{x}, \vec{\omega})L_e(\mathbf{x}, \vec{\omega})dz, \quad (2.1)$$

where $L(\mathbf{x}, \vec{\omega})$ is the radiance at point \mathbf{x} in direction $\vec{\omega}$, $L_e(\mathbf{x}, \vec{\omega})$ is the emitted radiance of the medium, and $\sigma(\mathbf{x}, \vec{\omega})$ is the (direction-dependent) attenuation coefficient.

2.1.2 Volume rendering equation

The solution to the RTE is given by the volume rendering equation:

$$L(\mathbf{x}, \vec{\omega}) = T(\mathbf{x}, \mathbf{x}_z)L(\mathbf{x}_z, \vec{\omega}) + \int_0^z T(\mathbf{x}, \mathbf{x}_t)\sigma(\mathbf{x}_t, \vec{\omega})L_e(\mathbf{x}_t, \vec{\omega})dt, \quad (2.2)$$

where $T(\mathbf{x}, \mathbf{x}_t)$ is the transmittance from \mathbf{x} to \mathbf{x}_t , given by:

$$T(\mathbf{x}, \mathbf{x}_t) = \exp\left(-\int_0^t \sigma(\mathbf{x}_s, \vec{\omega})ds\right). \quad (2.3)$$

Notably, we can also consider the probability distribution of the ray termination distance, or the *free-flight distribution*, whose probability density function is given by the product of the transmittance and the attenuation coefficient at the termination point:

$$p_{\mathbf{x}, \vec{\omega}}(z) = \sigma(\mathbf{x}_z, \vec{\omega})T(\mathbf{x}, \mathbf{x}_z). \quad (2.4)$$

Assuming no background light source and near and far bounds t_n and t_f , we can then write the expected color of a camera ray $\mathbf{r}(t) = \mathbf{o} + t\vec{\omega}$ as

$$C(\mathbf{r}) = \int_{t_n}^{t_f} \sigma(\mathbf{r}(t), \vec{\omega}) L_e(\mathbf{r}(t), \vec{\omega}) \exp\left(-\int_{t_n}^t \sigma(\mathbf{r}(s), \vec{\omega}) ds\right) dt \quad (2.5)$$

$$= \int_{t_n}^{t_f} p_{\mathbf{o}, \vec{\omega}}(t) L_e(\mathbf{r}(t), \vec{\omega}) dt \quad (2.6)$$

2.1.3 Discretization

In practice, as proposed by Max [15], when rendering a single ray, the volume rendering equation is often discretized by sampling the ray at regular intervals and approximating the integral with quadrature methods.

Given a ray $\mathbf{r}(t) = \mathbf{o} + t\vec{\omega}$ and discrete samples t_0, t_1, \dots, t_N along the ray, the color of the ray is approximated as

$$C(\mathbf{r}) \approx \sum_{i=0}^{N-1} p_i L_e(\mathbf{r}(t_i), \vec{\omega}), \quad (2.7)$$

where the free-flight distribution is also approximated as

$$p_i = \alpha_i \sum_{j=0}^{i-1} (1 - \alpha_j), \quad \alpha_i = 1 - \exp(-\sigma(\mathbf{r}(t_i), \vec{\omega}) \Delta t_i). \quad (2.8)$$

Notably, this reduces volume rendering to alpha compositing, where the color of the ray is accumulated by blending the color of each sample with the accumulated color.

Rendering a scene with the volume rendering equation requires knowing the attenuation coefficient σ and emitted radiance L_e at every point in the scene. Traditionally, these quantities are estimated using physical measurements and known material properties. However, more recently, neural networks have been used as a tool to directly estimate these quantities from images of the scene, as we will discuss in the next section.

2.2 Neural rendering

Recent works following the introduction of neural fields (NeRF) [17] have shown that neural networks are capable of representing complex scenes and rendering them with high fidelity. We briefly review the key concepts of neural rendering, which we will build upon in our work.

2.2.1 Neural fields

Neural fields are neural networks, or functions $f_{\Theta} : (\mathbf{x}, \vec{\omega}) \rightarrow (c, \sigma)$, that map a 3D spatial location $\mathbf{x} = (x, y, z)$ and a 2D viewing direction $\vec{\omega} = (\theta, \phi)$ to the emitted radiance and the attenuation coefficient at the given spatial location in the given viewing direction [17].

Neural fields are often trained to minimize the error between rendered pixels and ground truth pixels, which are obtained by sampling from a dataset of captured images of the scene. The

trained neural field can then be used to render novel views of the scene by sampling camera rays and evaluating the neural field at the desired spatial location and viewing directions.

Despite its success in rendering complex scenes, neural fields have limitations in extracting surfaces from the scene, as they do not explicitly model the geometry of the scene. Naively, one can also use neural fields to directly extract surfaces from the scene by thresholding the attenuation coefficient σ at a certain value. However, this approach is not ideal for extracting surfaces, and often results in incorrect or noisy surfaces.

To address this limitation, other works have proposed ways to directly model the geometry of the scene and connect the geometry of the scene to the attenuation coefficient for volume rendering, as we will discuss in the next section.

2.2.2 Neural surface representations

To address the limitations of using the attenuation coefficient to extract surfaces, works including [18, 21, 22] have proposed ways to directly model the *signed-distance function* (SDF) or *occupancy* of a scene and convert them into attenuation coefficients for volume rendering. These works have shown that neural surface representations can be used to both render high-quality images of scenes and extract surfaces from the scene, by using marching cubes [14] to find the 0.5-crossing of the occupancy, for example.

Specifically, given the occupancy of a scene, we adopt the method proposed by Miller et al. [18] and compute the attenuation coefficients by the formula

$$\sigma(\mathbf{x}, \vec{\omega}) = \frac{|\vec{\omega} \cdot \nabla o(\mathbf{x})|}{1 - o(\mathbf{x})}, \quad (2.9)$$

where $o : \mathbb{R}^3 \rightarrow [0, 1]$ is the occupancy function that represents the probability of a point in space being inside of an object. We further discuss how these concepts relate to our representation in section 3.1.

The original discrete formulation makes the assumption that the attenuation coefficient is constant within each discrete segment of the ray. As proposed by both Wang et al. [21] and Miller et al. [18], we can relax this assumption to allow for non-constant but monotone attenuation coefficients within each segment. Then, we can directly estimate α_i from the occupancy as

$$\alpha_i = \frac{|o(\mathbf{r}(t_i)) - o(\mathbf{r}(t_{i+1}))|}{1 - \min\{o(\mathbf{r}(t_i)), o(\mathbf{r}(t_{i+1}))\}}. \quad (2.10)$$

This alternative formulation also avoids the need to explicitly compute the attenuation coefficient and thus the gradient of the occupancy function, which requires additional computation and are often noisy.

Chapter 3

Winding number and dipole sums

3.1 The winding number

To lay the foundation for our point-based representation, we begin with an introduction on the winding number for surfaces and point clouds, which we will generalize in 3.2 for our implicit surface representation.

3.1.1 Winding number for surfaces

We first consider a continuous surface $\Gamma \subset \mathbb{R}^3$. There are many equivalent definitions of the winding number [6]; we follow Barill et al. [1] and use its definition as a *jump harmonic* scalar field. Then, the *winding number* $w : \mathbb{R}^3 \rightarrow \mathbb{R}$ is the scalar field solution to the Laplace boundary value problem (BVP) with jump Dirichlet and Neumann boundary conditions:

$$\Delta w(x) = 0 \text{ in } \mathbb{R}^3 \setminus \Gamma, \quad (3.1)$$

$$w^+(x) - w^-(x) = 1 \text{ on } \Gamma, \quad (3.2)$$

$$\partial w^+ / \partial n(x) - \partial w^- / \partial n(x) = 0 \text{ on } \Gamma. \quad (3.3)$$

Here, $n(x)$ is the normal at point $x \in \Gamma$, and $w^\pm(x) \equiv \lim_{\varepsilon \rightarrow 0} w(x \pm \varepsilon \cdot n(x))$ are the winding number values on either side of the surface Γ along the normal direction. Krutitskii [13] provide a detailed treatment of such BVPs, and in particular prove the following *boundary integral* expression for their solution:

$$w(x) = \int_{\Gamma} P(x, y) \cdot 1 \, d\sigma(y), \quad P(x, y) \equiv \frac{1}{4\pi} \frac{\langle n(y), \widehat{xy} \rangle}{\|x - y\|^2}, \quad (3.4)$$

where $\widehat{xy} \equiv (y - x) / \|y - x\|$ is the direction from x to y , and $P : \mathbb{R}^3 \times \mathbb{R}^3 \rightarrow \mathbb{R}$ is the *free-space Poisson kernel* for the Laplacian partial differential equation (PDE). We make explicit the factor 1 in the integral of 3.4, corresponding to the jump Dirichlet boundary condition 3.2, for reasons we will explain in 3.2.

When the surface Γ is the watertight boundary of one or more three-dimensional objects, then the winding number equals their binary *indicator function*— $w(x) = 1$ for points x at the objects' interior, $w(x) = 0$ otherwise.

3.1.2 Winding number for point clouds

We now consider an *oriented* point cloud $\mathcal{P} \equiv \{(p_m, n_m, A_m)\}_{m=1}^M$, where for each m we assume that:

1. the point p_m is a sample from an underlying surface Γ ;
2. the vector n_m is the normal of Γ at p_m ; and
3. the scalar A_m is the geodesic Voronoi area on Γ of p_m , i.e., the area of the subset of Γ where points are closer (in the geodesic distance sense) to p_m than any other point in the point cloud.

In practice, if only the points p_m are available, we can estimate the normals n_m and area weights A_m using standard techniques (e.g., by fitting and Voronoi-tessellating a plane to each point's k -nearest neighbors [8, 19]).

As Barill et al. [1] explain, the boundary integral representation 3.4 directly suggests the following generalization of the winding number for point clouds:

$$w_{\text{pc}}(x) \equiv \sum_{m=1}^M A_m P(x, p_m) \cdot 1 = \sum_{m=1}^M \frac{A_m}{4\pi} \frac{\langle n_m, \widehat{xp_m} \rangle}{\|x - p_m\|^2} \cdot 1. \quad (3.5)$$

Barill et al. [1] show that w_{pc} is a non-binary scalar field that approaches $1/2$ at points near the boundary of the continuous surface Γ underlying the point cloud \mathcal{P} , increases towards its interior, and decreases towards its exterior. Thus, the $1/2$ -level set of w_{pc} is an implicit surface that approximates Γ ; this approximation becomes exact as point density converges to infinity, and degrades gracefully as the number M of points decreases.

3.1.3 Barnes-Hut approximation

Evaluating the winding number $w_{\text{pc}}(x)$ at a query point x using 3.5 has linear complexity $O(M)$ with respect to the point cloud size M ; for large point clouds, doing so can be exceedingly expensive, especially if we need to query w_{pc} at multiple points (as we will later in this section). Barill et al. [1] show how to compute $w_{\text{pc}}(x)$ with logarithmic complexity using the *Barnes-Hut fast summation method* [2]. This method first creates a tree data structure (e.g., octree [16]) whose nodes hierarchically subdivide points in the point cloud into clusters, with leaf nodes corresponding to individual points. Each tree node t has a centroidal location, area-weighted normal, and radius

$$\tilde{p}_t \equiv \frac{\sum_{m \in \mathcal{L}(t)} A_m p_m}{\sum_{m \in \mathcal{L}(t)} A_m}, \quad \tilde{n}_t \equiv \sum_{m \in \mathcal{L}(t)} A_m n_m \cdot 1, \quad \tilde{r}_t \equiv \max_{m \in \mathcal{L}(t)} \|p_m - \tilde{p}_t\|. \quad (3.6)$$

where $\mathcal{L}(t)$ is the set of leaf (i.e., single-point) nodes that are successors of t in the tree hierarchy. We purposefully include an additional factor of 1 in the are-weighted normal computation, for reasons we will explain in 3.2.3.

Then, for each query point x , the Barnes-Hut methods performs a depth-first tree traversal; at each node t , if x is sufficiently far from its centroid (i.e., $\|x - \tilde{p}_t\| > \beta \tilde{r}_t$), the node's successors

are not visited and the sum of contributions to $w_{\text{pc}}(x)$ from all leaves in $\mathcal{L}(t)$ is approximated as:

$$\sum_{m \in \mathcal{L}(t)} A_m P(x, p_m) \cdot 1 \approx \tilde{P}(x, \tilde{p}_t) \equiv \frac{1}{4\pi} \frac{\langle \tilde{n}_t, \widehat{x\tilde{p}_t} \rangle}{\|x - \tilde{p}_t\|^2}. \quad (3.7)$$

This approximation expresses the fact that, due to the squared-distance falloff of the Poisson kernel in 3.4, the *far-field* influence of a cluster of points can be represented by a single point at the cluster’s centroid. We also note that the area weight is not included in the far-field approximation, as it is already accounted for in the normal computation in 3.6.

3.1.4 Rendering point clouds with winding numbers

Barill et al. [1] show that efficient winding number queries facilitate several point cloud operations, e.g., meshing, inside-outside tests, and Boolean composition. Our goal in this paper is to show that, with appropriate modifications (3.2), these queries facilitate also forward and inverse rendering of geometry represented as point clouds.

Specifically, we can use the winding number w_{pc} to define an implicit surface Γ_{pc}

$$\Gamma_{\text{pc}} \equiv \{x \in \mathbb{R}^3 : w_{\text{pc}}(x) = 1/2\}, \quad (3.8)$$

and, by naturally viewing the winding number as an occupancy function and following Miller et al. [18, Equation (12)], we can compute a volumetric attenuation coefficient σ_{pc} as:

$$\sigma_{\text{pc}}(x, \omega) \equiv \frac{|\omega \cdot \nabla w_{\text{pc}}(x)|}{1 - w_{\text{pc}}(x)} \quad (3.9)$$

Then, for surface rendering, we can perform ray casting and visibility queries on the point cloud, by intersecting the isosurface Γ_{pc} using ray marching [7]. Likewise, for volumetric rendering, we can compute free-flight distribution and transmittance queries through the point cloud, by accumulating the coefficient σ_{pc} along a ray. All these ray operations use only the point cloud attributes, and do not require meshing or using a proxy (e.g., grid or neural) for the point cloud. Additionally, though each ray operation requires winding number queries at multiple ray points, they remain efficient thanks to the Barnes-Hut method. Lastly, backpropagating through the expressions for σ_{pc} , and w_{pc} to update point cloud parameters is straightforward and efficient, as we discuss in 3.3.

Unfortunately, despite these attractive properties, the winding number w_{pc} —and associated isosurface Γ_{pc} and attenuation coefficient σ_{pc} —have a few critical shortcomings that make them unsuitable for direct use in rendering applications. We explain these shortcomings, and how to overcome them, in the next section.

3.1.5 Relationship to Poisson surface reconstruction

Before we continue, we remark on a relationship between the point cloud winding number w_{pc} and Poisson surface reconstruction [10, 11]. As Barill et al. [1] explain, both techniques compute, from an oriented point cloud, a scalar field that approximates the true winding number,

corresponding to the solution of BVP (3.1, 3.2, 3.3) for the continuous surface underlying the point cloud. The *limit* behaviors of the two scalar fields are equivalent. However, whereas the approximation of 3.5 can be efficiently computed directly from the point cloud, the approximation by Poisson surface reconstruction requires solving an expensive Poisson integration problem, making it impractical for forward and (especially) inverse rendering applications. Using w_{pc} allows us to efficiently render an approximation to the implicit surface output by Poisson surface reconstruction, without the need for a Poisson solver.

3.2 Fast dipole sums

We introduce a generalization of 3.5 that will serve as our point-based representation for both geometry and radiance in inverse rendering applications. We first derive our generalization, then explain its advantages.

3.2.1 General Dirichlet conditions

We generalize the BVP (3.1, 3.2, 3.3) to use an arbitrary *Dirichlet data* function $f : \Gamma \rightarrow \mathbb{R}$ for the jump Dirichlet boundary condition:

$$\Delta u(x) = 0 \text{ in } \mathbb{R}^3 \setminus \Gamma, \quad (3.10)$$

$$u^+(x) - u^-(x) = f(x) \text{ on } \Gamma, \quad (3.11)$$

$$\partial u^+/\partial n(x) - \partial u^-/\partial n(x) = 0 \text{ on } \Gamma. \quad (3.12)$$

We also augment the point cloud $\mathcal{P} := \{(p_m, n_m, A_m, f_m)\}_{m=1}^M$ to include the Dirichlet data as a per-point attribute, $f_m \equiv f(p_m)$. We will use u^f to denote the solution to this BVP for specific Dirichlet data f . Then, we can modify (3.4, 3.5) to express u^f and its point-based approximation as [13]:

$$u^f(x) \equiv \int_{\Gamma} P(x, y) \cdot f(y) d\sigma(y), \quad u_{pc}^f(x) \equiv \sum_{m=1}^M A_m P(x, p_m) \cdot f_m. \quad (3.13)$$

3.2.2 Regularized Poisson kernel

The Poisson kernel $P(x, y)$ is singular as $x \rightarrow y$; this makes the value of u_{pc}^f at locations x near a point p_m in the point cloud numerically unstable, and undefined at p_m .

To overcome this issue, we use the method of *regularized fundamental solutions* developed in PDE simulation [3, 4, 5] to address similar numerical issues from these singularities. Its starting point is the definition of the Poisson kernel through the *Green's function* (or *fundamental solution*) $G : \mathbb{R}^3 \times \mathbb{R}^3 \rightarrow \mathbb{R}$ of the Laplace PDE:

$$P(x, y) \equiv n(y) \cdot \nabla_x G(x, y) \text{ where } G \text{ satisfies } \Delta G(x, y) = \delta(x - y), \quad (3.14)$$

where δ is the Dirac delta distribution in \mathbb{R}^3 . The method of regularized fundamental solutions replaces δ with a *nascent delta function*, that is, a function $\phi_\varepsilon(x - y)$ satisfying $\lim_{\varepsilon \rightarrow 0} \phi_\varepsilon(x - y) =$

$\delta(x - y)$. Then, we can define the regularized Green's function G_ε and Poisson kernel P_ε exactly analogously to 3.14:

$$P_\varepsilon(x, y) \equiv n(y) \cdot \nabla_x G_\varepsilon(x, y) \text{ where } G_\varepsilon \text{ satisfies } \Delta G_\varepsilon(x, y) = \phi_\varepsilon(x - y), \quad (3.15)$$

where it follows that $\lim_{\varepsilon \rightarrow 0} G_\varepsilon = G$ and $\lim_{\varepsilon \rightarrow 0} P_\varepsilon = P$. A common nascent delta function is the Gaussian function $\phi_\varepsilon(x - y) \equiv 1/\varepsilon\sqrt{2\pi} \cdot \exp(-\|x - y\|^2/2\varepsilon^2)$. The corresponding regularized Poisson kernel equals [3]:

$$P_\varepsilon(x, y) \equiv P(x, y) \cdot S\left(\frac{\|x - y\|}{\varepsilon}\right), \text{ where } S(t) \equiv \text{erf}(t) - \frac{2}{\sqrt{\pi}} \cdot t \cdot \exp(-t^2). \quad (3.16)$$

The parameter ε provides control between regularization (restricting how fast P_ε increases as it approaches singularity) and bias (controlling the difference between P_ε and P).

3.2.3 Regularized dipole sum

As shown in 3.5 and 3.13, the point-cloud winding number is the solution to the BVP for constant unit data, i.e., $w_{\text{pc}}(x) = u_{\text{pc}}^1(x)$. However, as explained above, naively using the Poisson kernel for the dipole sum in 3.17 leads to numerical instability during both optimization and mesh extraction, as the kernel is singular near the surface. The regularized dipole sum in 3.17 overcomes this issue by using the regularized Poisson kernel, which is always finite and provides a stable approximation to the true dipole sum. Moreover, in practice, using constant unit data for the dipole sum is also not always ideal, as it is not robust to noisy point clouds and gaps and holes in the point cloud.

Therefore, we instead use the regularized dipole sum with non-unit Dirichlet data,

$$u_{\text{pc}, \varepsilon}^f(x) \equiv \sum_{m=1}^M A_m P_\varepsilon(x, p_m) \cdot f_m, \quad (3.17)$$

where ε is the regularization strength and f_m is the per-point Dirichlet data, both of which are learnable parameters. This allows us to accurately represent the underlying geometry of the point cloud, while also providing a stable and robust representation for rendering applications.

We note that although this dipole sum is written as a sum over all points in the point cloud, introducing the regularization term and non-unit Dirichlet data does not prevent us from using the Barnes-Hut method to efficiently compute the dipole sum with logarithmic complexity. Indeed, the regularized dipole sum can be computed using the same tree data structure but with a modified area-weighted normal that also takes into account the non-unit Dirichlet data:

$$\tilde{n}_t \equiv \sum_{m \in \mathcal{L}(t)} A_m n_m \cdot f_m \quad (3.18)$$

The contributions from the leaf nodes in the far-field approximation are approximated as:

$$\sum_{m \in \mathcal{L}(t)} A_m P_\varepsilon(x, p_m) \cdot f_m \approx \tilde{P}_\varepsilon(x, \tilde{p}_t) \equiv \frac{1}{4\pi} \frac{\langle \tilde{n}_t, \widehat{x\tilde{p}_t} \rangle}{\|x - \tilde{p}_t\|^2} S\left(\frac{\|x - \tilde{p}_t\|}{\varepsilon}\right). \quad (3.19)$$

Again, both the area weights and the Dirichlet data are already accounted for when computing the area-weighted normal in 3.18, and are not included in the far-field approximation in 3.19. Crucially, this is *not* the same as separately aggregating area weights and Dirichlet data for the node; instead, this correctly computes the summed contribution from the leaf nodes under the approximation that they are concentrated at the centroid of the node.

3.2.4 Rendering with regularized dipole sums

Finally, we introduce two modifications that makes the regularized dipole sum suitable for rendering applications.

First, although for watertight surfaces, the winding number is a binary indicator function, in practice, when discretized on a potentially noisy point cloud, the winding number can easily take on values greater than 1 or less than 0. To ensure that the occupancy function derived from the winding number is well-defined, we apply a *sigmoid* function to the regularized dipole sum:

$$o(x) \equiv \Phi_s \left(u_{\text{pc},\varepsilon}^f(x) - \frac{1}{2} \right), \quad \Phi_s(x) \equiv \frac{1}{1 + \exp(-sx)}, \quad (3.20)$$

where Φ_s is the logistic sigmoid with an additional scale parameter s , which maps the regularized dipole sum to the range $[0, 1]$.

The point cloud winding number does not guarantee a sharp transition between the inside and outside of the surface, which is crucial for representing sharp geometric features. The scale parameter of the sigmoid function allows us to control the sharpness of the transition without affecting the geometry of the underlying implicit surface, as the $1/2$ -level set remains unchanged.

Second, in addition to per-point Dirichlet data used to represent geometry, we can similarly store per-point neural features to also model the appearance of the scene. However, unlike for geometry, we do not want sharp discontinuity in the appearance of the scene near the surface. To address this, we use a modified version of the Poisson kernel to *interpolate* d -dimensional per-point neural features $h_m \in \mathbb{R}^d$ at query points x :

$$h(x) \equiv \sum_{m=1}^M A_m P^{\text{mod}}(x, p_m) \cdot h_m, \quad P^{\text{mod}}(x, y) \equiv \frac{1}{4\pi \|x - y\|^2}, \quad (3.21)$$

We omit the regularized version of the modified kernel and the corresponding Barnes-Hut approximation for brevity, as they are analogous to the original Poisson kernel. The modified kernel ensures that the neural features are smoothly interpolated across the scene without sharp discontinuities.

With interpolated neural features, we follow Wang et al. [21] and use a neural network to predict the radiance at a query point x with viewing direction ω and neural features $h(x)$:

$$L(x, \omega) \equiv N(x, \omega, h(x)), \quad (3.22)$$

In particular, we also follow Verbin et al. [20] and use spherical harmonics to encode the viewing direction instead of the positional encoding original proposed by Mildenhall et al. [17].

3.3 Logarithmic complexity backpropagation

In this section, we show how we can utilize the Barnes-Hut approximation to efficiently back-propagate gradients through the octree data structure, and how this allows us to efficiently optimize the point cloud parameters. For simplicity, we focus on deriving the gradient computation for the per-point Dirichlet data, but the same principles apply to the per-point neural features.

3.3.1 Gradient accumulation

Denoting the rendered pixels as \hat{C}_i and ground truth pixels as C_i , we can define the loss function as the mean absolute error between the rendered and ground truth pixels:

$$\text{loss} = \frac{1}{N} \sum_{i=1}^N \left\| \hat{C}_i - C_i \right\|_1. \quad (3.23)$$

Since we use the Barnes-Hut approximation during forward rendering, the loss can be viewed as a function of both the per-point parameters of the point cloud and per-node parameters of the octree structure.

Assuming we have a point cloud of n points and make m winding number queries during rendering, naively using autograd to compute the gradients of the loss with respect to all of the parameters would require $O(n \cdot m)$ time at best: for each single query, backpropagation requires computing the gradient at all octree nodes that are visited during forward rendering. This is identical to what happens during forward rendering and only takes $O(\log n)$ time. However, since the per-node parameters are essentially functions of the per-point parameters of its leaves, this means that for each node t we visit, we would also need to visit all of its leaves $\mathcal{L}(t)$ and accumulate gradients at its leaves. This results in asymptotically slower backpropagation compared to forward rendering.

To address this, we introduce a two-stage backpropagation scheme:

1. In the first stage, we *detach* the per-node parameters from the per-point parameters and compute the gradients of the loss with respect to only the per-node parameters. For m winding number queries during rendering, this only requires $O(m \log n)$ time, identical to forward rendering.
2. In the second stage, given the gradients of the per-node parameters, we can use the chain rule to compute the gradient of the per-point parameters. Since each point can only be the leaf of $O(\log n)$ many nodes, in the worst case, this step still only requires $O(n \log n)$ time. Importantly, this is only a one-time cost for each iteration of training, and does not need to be done for each individual query.

In summary, this two stage backpropagation scheme allows us to run backpropagation in $O((n + m) \log n)$ time, compared to the naive $O(n \cdot m)$ time. In practice, n and m can be on the order of thousands and millions, respectively, making this optimization crucial for efficient training.

3.3.2 Gradient computation for nodes

Next, we elaborate on how the gradient of the loss with respect to the per-node parameters can be computed efficiently, which is essential to the first stage of our backpropagation scheme. For each point m , we denote its Dirichlet data as f_m and the gradient of the per-point parameters as ∇f_m .

For each node t , we note that the per-node Dirichlet data is not explicitly stored but as part of the area-weighted normal \tilde{n}_t as shown in 3.18. In other words, instead of computing the gradient of the per-node Dirichlet data, we need to instead compute the gradient of area-weighted normal, which we denote as $\nabla \tilde{n}_t$.

We consider a single node t and all queries $Q(t)$ that visit this node during forward rendering. For each query, we denote the corresponding regularized dipole sum as $u_{pc,\varepsilon}^f(x)$ and the gradient of the regularized dipole sum as $\nabla u_{pc,\varepsilon}^f(x)$. We can then compute the gradient of the area-weighted normal as:

$$\nabla \tilde{n}_t = \sum_{x \in Q(t)} \nabla u_{pc,\varepsilon}^f(x) \cdot \frac{1}{4\pi} \frac{\widehat{x\tilde{p}_t}}{\|x - \tilde{p}_t\|^2} S\left(\frac{\|x - \tilde{p}_t\|}{\varepsilon}\right). \quad (3.24)$$

In practice, we loop over all queries and traverse the octree to accumulate the gradients of the area-weighted normal at each node. This allows us to compute the gradient of the loss with respect to the per-node parameters in $O(m \log n)$ time.

3.3.3 Propagating gradient to points

Then, the second stage of backpropagation requires propagating gradients stored on the nodes of the octree down to its leaves, i.e., the individual points of the point cloud. During the construction of the octree structure, we computed the area-weighted normal for each node as shown in 3.18. Given the gradient of the area-weighted normal $\nabla \tilde{n}_t$, we can compute the gradient of the per-point Dirichlet data as:

$$\nabla f_m = \sum_{t \in \mathcal{T}(m)} \langle \nabla \tilde{n}_t \cdot A_m n_m \rangle, \quad (3.25)$$

where $\mathcal{T}(m)$ is the set of nodes that are ancestors of point m in the octree. In practice, we loop over all nodes of the octree and propagate gradients down to the leaves of each node.

The second stage can be done independently of the first stage, and only needs to be done once for each iteration of training. This allows us to compute the gradient of the loss with respect to the per-point parameters in an additional $O(n \log n)$ time.

3.3.4 Optimization

Finally, we can use the gradients computed in the two stages to optimize the per-point parameters of the point cloud. We can use any optimization algorithm, such as Adam [12], to update the per-point parameters.

After updating the per-point parameters, we can then recompute the new area-weighted normal for each node of the octree efficiently in $O(n \log n)$ time, according to the same equation as in 3.18. This summarizes how we can efficiently optimize the learnable parameters of the point cloud using the Barnes-Hut approximation in a differentiable rendering setting.

Chapter 4

Experimental evaluation

4.1 3D reconstruction

4.1.1 Implementation details

4.1.2 Comparisons

4.2 Multi-bounce rendering

4.2.1 Experimental setup

4.2.2 Results

4.3 Visualizations of various quantities

4.3.1 Fields

4.3.2 Point clouds

Chapter 5

Conclusion

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