Multi-view Stereo Reconstruction of Dense Shape and Complex

Appearance

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Abstract

We address the problem of estimating the three-dimensional shape and complex appearance of a scene from a calibrated set of views under fixed illumination. Our approach relies on an affine subspace constraint that must be satisfied when the scene exhibits "diffuse + specular" reflectance characteristics. This constraint is used to define a cost functional for the discrepancy between the measured images and those generated by the estimate of the scene, rather than attempting to match image-to-image directly. Minimizing such a functional yields the optimal estimate of the shape of the scene, represented by a dense surface, as well as its radiance, represented by three functions defined on such a surface. These can be used to generate novel views that capture the non-Lambertian appearance of the scene.

1 Introduction

Multi-frame stereo consists of reconstructing the three-dimensional (3-D) shape of a scene from a collection of images taken from different vantage points. This is one of the classical problems of computer vision, where significant progress has been made in the last decade. In the early days of stereo, it was common to decompose the problem into two steps: establishing correspondence between points in different views, and then triangulating their position in space. Points in different images are said to be in correspondence when they are images of the same physical point in space via perspective projection. Once correspondence is established, the position of the points as well as the relative pose of the cameras can be determined using well-established procedures that are now the subject of textbooks [7, 11, 22].

Unfortunately the first step, establishing correspondence, is far less amenable to a clean and simple solution. First of all, point correspondence can only be reliably established for a very small subset of the

scene. For instance, given a scene that contains a white wall, we cannot say which point on one image of the wall corresponds to in another image, since the local appearance is the same for every neighborhood of a point. Therefore, after establishing correspondence and reconstructing the 3-D position of relatively few "feature" points,¹ one would have to "densify" the reconstruction by filling in points that cannot be matched from image to image.

Second, and more important, correspondence cannot be established by just comparing local image statistics unless the scene has the property that its appearance does not change with the viewpoint. Materials that exhibit this property are called Lambertian, or diffuse, and they include matte surfaces such as chalk, rough stone and certain fabrics². However, most of the materials that populate our daily scenes such as plastic, polished stone, skin, glass, metal, etc. do not enjoy this property. Indeed, one can make an object that deviates severely from the ideal Lambertian model, for instance a car, appear arbitrarily different from image to image by changing the viewpoint and the illumination.

In this paper the first issue is addressed at the outset by modeling the shape of the scene as a collection of smooth surfaces: Like many recent works in multi-view stereo, we do not seek to establish correspondence among a sparse set of feature points and then fill in the rest. Rather, we start with a generic surface, say a large sphere or a smooth cube, and evolve it, possibly via changes of topology, to best approximate the shape of the scene. We do so by numerically integrating systems of partial differential equations using the level set method. The second issue is addressed by bypassing the direct comparison of local image intensity, and instead comparing all images to the underlying model of the scene, which necessarily includes the current estimate of its shape *as well as its radiance*. Our model of the radiance is

¹Even a few thousand feature points are far fewer than the millions of pixels in an image

²Most feature correspondence algorithms implicitly assume that the scene is "almost" Lambertian, in the sense that the deviation from an ideal Lambertian model is small, not modeled explicitly, and instead lumped together with other factors as "noise."

not in an explicit functional form; instead, it accounts for deviations from Lambertian reflection through an affine subspace constraint on the radiance tensor field, which we will define shortly³.

The result is an algorithm that takes as input a sequence of images of a scene with complex appearance, such as those in Figure 1 and, with no intermediate steps, returns an estimate of its shape, described by one or more "dense" surfaces, and an estimate of its appearance, described by the radiance tensor field. Such a description can be used to render the scene from novel viewpoints, assuming a static illumination, in ways that preserve the complex appearance of the original scene.

Since a general scene cannot be reconstructed under varying and unknown illumination, we must make assumptions about the imaging process. Specifically, we assume that illumination is fixed but otherwise arbitrary, except for being "far enough" from the scene in a way that we will make precise in Section 2.1. Furthermore, we assume that the scene is a collection of smooth or piecewise smooth surfaces, and that its reflectance can be modeled by the linear combination of an ideal Lambertian component and a specular component, or what is known in computer graphics as a "diffuse + specular" reflectance model.

In the next subsection we will briefly review the state of the art as it relates to our contribution. Before we formally introduce the quantities at play in Section 2.1 we use the terms "photometry," "radiance," "reflectance" and "appearance" interchangeably, and similarly for "shape," "structure" and "geometry."

³Incidentally, as shown in [28], the distinction of comparing all images to an underlying model, as opposed to matching image-to-image, is relevant only in the presence of non-Lambertian scenes, or other constraints on the diffuse albedo, as we will discuss shortly and as shown in [28].

1.1 Relation to prior work

In order for any 3-D reconstruction to be possible, some assumption must be made on the photometry of the scene⁴. The most common assumption is that the light is fixed and the scene is *Lambertian*, i.e., the energy radiated from any point in the scene does not depend on the outgoing direction, so that correspondence can be easily established by comparing individual images. Indeed, as we have shown in [28], under the assumption of Lambertian reflection and in the absence of any additional information or constraint on the diffuse albedo, there is no difference between comparing all images to an underlying model of the scene as opposed to matching image-to-image directly. The situation is quite different, as we discuss in [28], when the scene deviates from ideal Lambertian reflection. In this case, reflection is most often described by an explicit model, a bidirectional reflectance distribution function (BRDF), chosen among a parametric class derived by physical or empirical considerations⁵.

Most often, however, deviations from Lambertian reflection are modeled as "noise" or "outliers" and either minimized by choice of a suitable cost functionals (such as photo-consistency [21]), or rejected using robust statistical methods. For instance, one can select candidates for correspondence in each image by looking at image statistics integrated over a region, compute the cross-correlation or other score among putative correspondences, and then test whether they are consistent with a common epipolar geometry. This works well when the scene is composed mostly of matte surfaces with few specular highlights. However, for objects that are *shiny* and concentrated light distributions (see Figure1), this approach shows limitations. Alternatively, one can set up a global cost functional obtained by integrating

⁴It is straightforward to show that if a scene has arbitrary reflectance properties and one can change the light distribution from frame to frame, correspondence can*not* be established [32].

⁵For instance, [23] and [33] exploit the reciprocity condition of the BRDF to perform reconstruction using a particular imaging setup where multiple images are obtained by swapping a point light source and the camera. We do not impose constraints on the viewpoint, and do not restrict the illumination to be a point source. Indeed, we do not model illumination explicitly, so our approach is quite different.



Figure 1: (COLOR) Scenes with strong specularities are a challenge to algorithms relying on image-toimage matching.

on the entire scene a local consistency measure (e.g. normalized cross-correlation) computed on the images, and minimize it with respect to the unknown shape using variational techniques, an approach pioneered by [8] in stereo reconstruction. Our approach is based on a similar philosophy, and we also uses level set methods [27] to numerically solve the variational problem. However, while Faugeras and Keriven estimate geometry alone, we estimate both geometry and photometry (radiance) and forego the Lambertian assumption that is latent in the cost functional used in [8]. [16] have modified the cost functional to minimize the effects of isolated specularities.

This work also relates to a series of works where the same computational framework is used in estimating the shape and radiance for scenes of increasing complexity: from constant diffuse albedo [31] to smooth diffuse albedo [18], to piecewise constant and piecewise smooth diffuse albedo [17], to arbitrary diffuse and constant specular albedo, to arbitrary diffuse and specular albedos.

In addressing non-Lambertian reflection, our work relates to several studies on specular reflections in stereo matching and reconstruction. [1] consider the likelihood of correct stereo matching by analyzing

the relationship between stereo vergence and surface roughness, and propose a trinocular system where only two images are used at a time in the computation of depth at a point. [2, 3] excise specularities as a pre-processing step, similarly to [26], while [24] do so using polarized filters.

Non-Lambertian reflection has also been addressed in the context of photometric stereo, for instance by [14]. In this case, the viewpoint is fixed while the illumination changes. In this respect, this is quite different from our approach, that is more in line with traditional multi-view stereo in assuming that the viewpoint, and not the illumination, moves. Other approaches [12] compare the observed images with that of objects with known shape to obtain surface normals and hence shape.

This work also relates to the general problem of estimating reflectance properties as well as shape from sequences of images. For instance [32] use known shape to estimate global illumination [32]; in light field rendering [5, 10, 25] there is no explicit reconstruction of shape, and the radiance tensor, extended to the volume, is sampled directly. Indeed, the rank constraint on the radiance tensor field is often used in light field rendering, albeit not for inferring properties of the scene but, rather, for computational efficiency.

This work addresses the problem of multi-view stereo with fixed illumination and arbitrarily changing viewpoint. To the best of our knowledge, we are the first to propose a multi-view stereo algorithm that can provide an estimate of both dense shape and non-Lambertian reflection. Our algorithm is based on an *affine subspace constraint on* the radiance tensor field (Section 2.1), which we show to imply (and hence be more general than) a "diffuse + specular" reflection models commonly used in Computer Graphics (Proposition 1).

2 Local modeling of radiance and image discrepancy

In this section we introduce the model of photometry, based on the *radiance tensor field*, and the measure of discrepancy between *model and images* that is the basis of our approach.

2.1 The radiance tensor field

Let S be a (smooth) surface embedded in \mathbb{R}^3 and P be the generic point on S, with coordinates $\mathbf{X} = [X_1, X_2, X_3] \in \mathbb{R}^3$ with respect to a fixed world reference frame. We denote with T_PS the tangent plane to the surface at the point P. The generic vector in the tangent plane (embedded in Euclidean space) has coordinates $v \in \mathbb{R}^3$. Let an ideal perspective camera be characterized by a Euclidean reference frame $g \in SE(3)$, that describes the change of coordinates between the world reference frame and the frame attached to the optical center of the camera, represented by a rotation matrix and a translation vector⁶. Therefore, if $\pi : \mathbb{R}^3 \longrightarrow \mathbb{R}^2$ denotes the canonical perspective projection⁷, the point P projects onto each image in the coordinates $\mathbf{x} = \pi(gP)$.

For each point $P \in S$ we consider a discretization of a small neighborhood $\Omega_P \subset T_PS$ around it. This discretization is usually done with a tessellation of T_PS , which we represent via the vectors v_1, v_2, \ldots, v_m , where m is the number of points in Ω_P , as shown in Figure 2. We assume to be able to measure the amount of light leaving these points toward a discrete number n of camera poses, g_1, g_2, \ldots, g_n . Therefore, to each point P we can associate an array of $m \times n$ ideal measurements, one

⁶ g acts on a point P with coordinates **X** via gP, which has coordinates **RX** + **T** where **R** \in SO(3) is an orthonormal matrix with positive determinant and **T** $\in \mathbb{R}^3$. The push-forward action of g on vectors $v \in T\mathbb{R}^3$ with coordinates **V** is given by g_*v , which has coordinates **RV**.

 $^{^{7}\}pi(\mathbf{X}) = [X_1/X_3, X_2/X_3].$



Figure 2: The local coordinate frame on the tangent plane, the discretization of the local neighborhood, and the projection onto an image.

column for each camera view and one row for each point in Ω_P , as

$$R(P) = \begin{bmatrix} \rho(v_1, g_1) & \dots & \rho(v_1, g_n) \\ \vdots & \ddots & \vdots \\ \rho(v_m, g_1) & \dots & \rho(v_m, g_n) \end{bmatrix}$$
(1)

where $\rho(v_i, g_j)$ can be thought of as an approximate measurement of the radiance of the surface at a point. Notice that $R_{ij} \doteq \rho(v_i, g_j)$ relates to the *ideal* image I_j with an explicit dependence on P via the irradiance equation [13, page 208], assuming a pin-hole projection:

$$R_{ij} = I_j(\pi(g_j(P+v_i))) \quad \forall v_i \in \Omega_P$$
⁽²⁾

for all j = 1, 2, ..., n. The map $S \to \mathbb{R}^{m \times n}$; $P \mapsto R(P)$ thus defines a tensor field on $S, R(\cdot)$ which, for any fixed P, is an $m \times n$ matrix, called the *radiance tensor*, or simply "radiance". In practice, the images I_j are measured only up to noise, so what is available is

$$\tilde{I}_j(\mathbf{x}) = I_j(\mathbf{x}) + w_j(\mathbf{x}); \quad \tilde{R}_{ij} = R_{ij} + w_{ij}$$
(3)

where $w_j(\mathbf{x})$ measures the discrepancy of the data from the model and can be considered as the realization of a random process (and therefore assumed to have a distribution associated to it), or simply as an unknown matrix whose norm we wish to minimize. We call \tilde{R} the measured radiance tensor field obtained by substituting the noisy images \tilde{I} in equation (2).

In general, the radiance tensor depends on the material properties of the surface and the lighting conditions. For instance, for the simplest case of Lambertian reflection,

$$R(P) = R_1(P) \cdot \mathbf{1}_n^T \tag{4}$$

where $R_j(P)$ denotes the *j*-th column of R(P) and $\mathbf{1}_n$ denotes an *n*-dimensional vector with all the elements equal to 1. It is because, by the Lambertian assumption, the radiance is independent of the viewpoint, and therefore all the columns of R(P) are identical. In fact, we can replace $R_1(P)$ in equation (4) with any other column of R(P). For more complex materials, R(P) has more structures but is, in general, not arbitrary. Proposition 1 shows that for ideal surfaces that obey a "diffuse+specular" reflection model, all the columns of R(R) live in an affine subspace of dimension 1. In order to set up the notation to state the proposition, we choose a reference frame $\langle \mathbf{e}_1, \mathbf{e}_2 \rangle$ for the tangent plane T_PS with the origin at $P:\langle \mathbf{e}_1, \mathbf{e}_1 \rangle = 1$, $\langle \mathbf{e}_2, \mathbf{e}_2 \rangle = 1$, $\langle \mathbf{e}_1, \mathbf{e}_2 \rangle = 0$. Let N_PS be the outward unit normal to S at



Figure 3: Illustration of the light interaction with the surface. H_P^2 is the unit hemisphere at a point P. $\lambda_i \in H_P^2$ is the incoming light direction and $\lambda_o \in H_P^2$ is the outgoing light direction. $\mathbf{h} \in H_P^2$ is the halfway vector between λ_i and λ_o . δ is the angle between \mathbf{h} and N. Surface reflectance is described by the bidirectional reflectance distribution function β that measures the ratio between the reflected energy along the direction λ_o due to the energy coming from the direction λ_i and the incoming energy.

P, so that $\mathbf{e}_1 \times \mathbf{e}_2 = N_P S$. Then $\langle \mathbf{e}_1, \mathbf{e}_2, N \rangle$ forms a Euclidean reference frame for \mathbb{R}^3 around *P*, where we have indicated the normal vector with *N* as a short-hand for $N_P S$. We denote with g_P the change of coordinates between the world reference frame and $\langle \mathbf{e}_1, \mathbf{e}_2, N \rangle$ (see Figure 2). We can parameterize each unit vector λ in the upper hemisphere at *P*, H_P^2 , with polar coordinates ($\theta_{\lambda}, \phi_{\lambda}$) $\in [0, \pi/2] \times [0, 2\pi]$, i.e., θ_{λ} is the angle between λ and *N* and ϕ_{λ} is the angle between λ and \mathbf{e}_1 , for all $\lambda \in H_P^2$.

The interaction of light with the surface S can be expressed, for most materials that we are going to deal with, by the *bidirectional reflectance distribution function* (BRDF⁸). This is a function of two directions in H_P^2 , the incident direction λ_i , parameterized by (θ_i, ϕ_i) and the reflected direction λ_o , parameterized by (θ_o, ϕ_o) , as well as the wavelength and polarization of the incident radiation, which we will ignore (see Figure 3). Ward's (anisotropic) elliptical Gaussian model [30] approximates the BRDF

⁸The BRDF is a simplified description of the radiometry of purely reflective (ideal) materials that yields an approximation of the radiance commonly used in computer graphics. It measures the ratio between the reflected energy along the direction (θ_o, ϕ_o) due to the energy coming from the direction (θ_i, ϕ_i) and the incoming energy.

 β with a linear combination of a diffuse term and a specular term:

$$\beta(\theta_i, \phi_i, \theta_o, \phi_o) = \frac{\rho_d}{\pi} + \frac{\rho_s \exp(-\tan^2 \delta(\cos^2 \gamma / \alpha_x^2 + \sin^2 \gamma / \alpha_y^2))}{4\pi \alpha_x \alpha_y \sqrt{\cos \theta_i \cos \theta_o}}$$
(5)

where ρ_d is the diffuse reflectance coefficient (commonly referred to as the *albedo*) and ρ_s is the specular reflectance coefficient; α_x and α_y are the standard deviations of the microscopic surface slope (surface roughness) in the direction of \mathbf{e}_1 and \mathbf{e}_2 respectively. They roughness coefficients are related to the properties of the material and we will consider them to be constant in a neighborhood of P. We note that constant surface roughness in a neighborhood does not imply that either diffuse reflectance coefficient (albedo) or specular reflectance coefficient is constant in that neighborhood. Let \mathbf{h} be the halfway vector between the directions λ_i and λ_o : $\mathbf{h} \doteq \frac{\lambda_i + \lambda_o}{\|\lambda_i + \lambda_o\|}$; (δ, γ) are the polar coordinates for \mathbf{h} and are therefore functions of $(\theta_i, \phi_i, \theta_o, \phi_o)$. The radiance in the direction determined by the point $\mathbf{x}_j = \pi(g_j P)$ in the *j*-th camera view is given by integrating the BRDF against the light distribution L in all directions (θ_i, ϕ_i) :

$$\rho(0,g_j) = \int_0^{2\pi} \int_0^{\pi/2} \beta(\theta_i,\phi_i,\theta_o,\phi_o) L(\theta_i,\phi_i) \cos\theta_i \sin\theta_i d\theta_i d\phi_i$$
(6)

where the direction from P to c_j , the *j*-th camera center, in the frame of the point P, i.e., $g_{P_*}^{-1}\left(\frac{c_j-P}{\|c_j-P\|}\right)$ (see Footnote 6), is represented in polar coordinates as (θ_o, ϕ_o) .

Proposition 1 (radiance tensor rank). Let S be made of a material that obeys the reflectance model (5). Furthermore, consider a surface patch $\Omega_P \subset T_P S$ that is small compared to the distance of P from the light sources and from the cameras. Then, if R(P) is computed for $v_i \in \Omega_P$ as in equation (1), we have that $\forall P \in S$

$$R_j(P) \in \mathbb{A}_m^1, \quad j = 1, 2, \dots, n \tag{7}$$

where \mathbb{A}_m^1 with $1 \leq r \leq \min(m, n) - 1$ denotes a r-dimensional affine subspace of \mathbb{R}^m .

Proof. To facilitate computing the radiance $\rho(v_i, g_j)$ for each $v_i \in \Omega_P \subset T_PS$; i = 1, 2, ..., m, in the direction of the origin of the reference frame of camera j = 1, 2, ..., n, we will denote with $\tilde{g}_j(v_i)$ the direction $g_{(P+v_i)*}^{-1}\left(\frac{c_j-(P+v_i)}{\|c_j-(P+v_i)\|}\right)$ from $P + v_i$ to c_j in the frame at the point $P + v_i : \langle \mathbf{e}_1(v_i), \mathbf{e}_2(v_i), N \rangle$. Since T_PS is a plane, we can choose $\langle \mathbf{e}_1(v_i), \mathbf{e}_2(v_i), N \rangle$ to coincide with the reference frame at $P : \langle \mathbf{e}_1, \mathbf{e}_2, N \rangle$. Under the assumption that Ω_P is small, we can approximate $\tilde{g}_j(v_i)$ with $\tilde{g}_j(0)^9$. Again, (θ_o, ϕ_o) are the polar coordinates of $\tilde{g}_j(0)$. Under the same assumption, we can also approximate the incoming light distribution at the point $P + v_i$ with $L(\theta_i, \phi_i)$. If we denote with $\rho(v_i|w)$ the radiance of point v_i along the direction λ , by equation (6), the radiance in the direction toward c_j is given by

$$\rho(v_i, g_j) = \rho\left(v_i | \tilde{g}_j(v_i)\right) \cong \rho\left(v_i | \tilde{g}_j(0)\right)$$

$$= \int \beta(v_i, \theta_i, \phi_i, \theta_o, \phi_o) L(\theta_i, \phi_i) \cos \theta_i \sin \theta_i d\theta_i d\phi_i$$

$$= \int_0^{2\pi} \int_0^{\pi/2} \frac{\rho_d(v_i)}{\pi} L(\theta_i, \phi_i) \cos \theta_i \sin \theta_i d\theta_i d\phi_i$$

$$+ \int \frac{\rho_s(v_i) \exp(-\tan^2 \delta(\cos^2 \gamma / \alpha_x^2 + \sin^2 \gamma / \alpha_y^2))}{4\pi \alpha_x \alpha_y \sqrt{\cos \theta_i \cos \theta_o}} L(\theta_i, \phi_i) \cos \theta_i \sin \theta_i d\theta_i d\phi_i$$

$$= \rho_d(v_i) s_0 + \rho_s(v_i) s_1(g_j)$$

where

$$s_{0} \doteq \int_{0}^{2\pi} \int_{0}^{\pi/2} \frac{1}{\pi} L(\theta_{i}, \phi_{i}) \cos \theta_{i} \sin \theta_{i} d\theta_{i} d\phi_{i}$$

$$s_{1}(g_{j}) \doteq s_{1}(\theta_{o}, \phi_{o}) = \int_{0}^{2\pi} \int_{0}^{\pi/2} \frac{\exp(-\tan^{2} \delta(\cos^{2} \gamma/\alpha_{x}^{2} + \sin^{2} \gamma/\alpha_{y}^{2}))}{4\pi \alpha_{x} \alpha_{y} \sqrt{\cos \theta_{i} \cos \theta_{o}}} L(\theta_{i}, \phi_{i}) \cos \theta_{i} \sin \theta_{i} d\theta_{i} d\phi_{i}.$$

⁹The meaning of approximation goes as follows: $\forall \epsilon > 0$, we can choose the size of Ω_P small such that $\|\tilde{g}_j(v_i) - \tilde{g}_j(0)\| < \epsilon$.

The fact that s_0 does not depend on g_j implies $R_j(P) \doteq [\rho(v_1, g_j), \rho(v_2, g_j), \dots, \rho(v_m, g_j)]^T \in \mathbb{A}_m^1$. \Box

Remark 1. ρ_d , s_0 , ρ_s and s_1 are all functions of P. The affine subspace \mathbb{A}^1_m may change with respect to P as well. We do not assume at all that either the albedos or the light distribution is the same for every P, but only that the surface roughness is locally constant in Ω_P .

Remark 2. Using $\tilde{g}_j(v_i)$ to approximate $\tilde{g}_j(v_0)$ is equivalent to using a scaled orthographic projection for the imaging model in Ω_P . However, when P moves over the surface, the parameters for the scaled orthographic projection are allowed to change. Therefore, we are not enforcing a scaled orthographic projection for the entire scene. The imaging model is still the perspective projection we put up at the beginning. In other words, we do not assume the overall size of the scene is small with respect to the distances to light sources or cameras.

Remark 3. Proposition 1 is stronger than that of [15]. The affine subspace constraint implies the rank constraint of [15], but the converse is not true. We will show in Section 3.1 that Proposition 1 leads to a simpler and faster algorithm than the one presented in [15].

The intuition behind this proposition is that, in the limit where the light sources are far, and the patch Ω_P is small, the specular component of the radiance of Ω_P is modulated by a scalar function that depends on the viewpoint ¹⁰. Of course, these conditions are a mathematical idealization. In practice, we verify experimentally that the dimension of \mathbb{A}_m^r decreases sharply and is negligible beyond the first. In the experimental section (Section 4) we will report how the sizes of neighborhoods affect the performance of the algorithm and we will also discuss the range of applicability of this constraint on realistic imaging conditions.

¹⁰Also note that the limit where the area of Ω_P goes to zero does not cause the rank to go to zero because the matrix R(P) becomes smaller, since one can resize the tessellation of the tangent plane so as to keep the number of rows of R(P) constant.

Regardless of the actual dimension for the affine subspace, a limitation on the dimension can be exploited to set up a discrepancy function for stereo reconstruction, as we do in the next section. In view of the claim above, one can then express the radiance tensor in terms of three vectors by lumping s_0 into d_0 . The relevance of Proposition 1 will be further discussed in Section 5.

Corollary 1 (local radiance model). At each point P of an ideal surface S that obeys the conditions of Proposition 1, the radiance tensor field can be represented with three vectors $d_0(v), d_1(v) \in \mathbb{R}^m$ and $s_1(g) \in \mathbb{R}^n$ as:

$$R(P) = d_0(v)\mathbf{1}_n^T + d_1(v)s_1^T(g).$$
(8)

As we have pointed in Remark 1, the reader should notice that $d_0(v)$, $d_1(v)$ and $s_1(g)$ are all functions of the point P on the surface. Note that s_1 depends on the viewing directions, a necessary element in modeling non-Lambertian reflection.

2.2 A discrepancy measure for non-Lambertian scenes

Naturally, due to image noise and deviation from the "diffuse+specular" reflectance model, the *measured* tensor $\tilde{R}(P)$ has its columns live in a higher dimensional affine subspace, or even \mathbb{R}^m . The key idea here is to use this subspace constraint to set up a matching criterion for stereo reconstruction. This is done by setting up an error function between the measured radiance tensor $\tilde{R}(P)$ and the model R(P) at each point P (see equation (3)):

$$\Phi(P) \doteq \|\tilde{R}(P) - d_0(v)\mathbf{1}_n^T - d_1(v)s_1^T(g)\|_F^2$$
(9)

where we have chosen the squared Frobenius norm to compare radiance tensors. Clearly $\Phi(P)$ will

depend on the coordinates of P. In addition, $\Phi(P)$ will also depend on the normal at P, since v_i lives in $T_PS: \Phi(P) = \Phi(\mathbf{X}, N)$. If we define

$$\phi_{ij} = \tilde{R}_{ij} - d_0(v_i) \mathbf{1}_n^T - d_1(v_i) s_1(g_j), \tag{10}$$

where \tilde{R}_{ij} is the (i, j)-th element of $\tilde{R}(P)$, $d_k(v_i)$ is the *i*-th component of $d_k(v)$ for k = 1, 2 and $s_1(g_j)$ is the *j*-th component of $s_1(g)$, then the squared Frobenius norm is the sum of the square of each element ϕ_{ij} . The surface *S* can then be found as the minimizer of the energy $E \doteq \int_S \Phi(P) dA$:

$$\hat{S} \doteq \arg\min_{S} \int_{S} \Phi(P) dA \tag{11}$$

where dA is the area measure on S.

For each *P*, we first compute the mean column of $\tilde{R}(P)$

$$\tilde{d}_0(v) \doteq \frac{1}{n} \sum_{j=1}^n \tilde{R}_j(P) \tag{12}$$

and then obtain the residual matrix $\widehat{R}(P)$ which is defined as

$$\widehat{R}(P) \doteq \widetilde{R}(P) - \widetilde{d}_0(v) \mathbf{1}_n^T.$$
(13)

As we have noted, since the actual measured tensor \tilde{R} will in general have no subspace constraint, the residual matrix $\hat{R}(P)$ will not have either. We can write it using the singular value decomposition (SVD)

as

$$\widehat{R}(P) = \sum_{i=1}^{r} \widetilde{d}_i(v) \widetilde{s}_i^T(g)$$
(14)

where r is the rank of $\widehat{R}(P)^{11}$. The singular values are sorted in a decreasing order with respect to k. Since, from the subspace constraint of Proposition 1, we can choose the basis of R arbitrarily, we can have

$$d_0(v) = \tilde{d}_0(v), \quad d_1(v) = \tilde{d}_1(v) \quad \text{and} \quad s_1(g) = \tilde{s}_1(g)$$
 (15)

and $R(P) = \tilde{d}_0(v)\tilde{\mathbf{1}}_n^T(g) + \tilde{d}_1(v)\tilde{s}_1^T(g)$. The function Φ can therefore be written as

$$\Phi(P) = \|\tilde{d}_2(v)\tilde{s}_2^T(g) + \tilde{d}_3(v)\tilde{s}_3^T(g) + \dots + \tilde{d}_r(v)\tilde{s}_r^T(g)\|_F^2.$$
(16)

Since $\tilde{d}_0(v)$ is the mean column of $\tilde{R}(P)$, we have that the columns of the residual matrix $\hat{R}(P)$ sum to zero, i.e.,

$$\sum_{j=1}^{n} \widehat{R}_{j}(P) = \sum_{j=1}^{n} \sum_{k=1}^{r} \widetilde{d}_{k}(v) \widetilde{s}_{k}^{T}(g_{j}) = 0$$
(17)

where $\tilde{s}_k(g_j)$ is the *j*-th component of $\tilde{s}_k(g)$. By the properties of the SVD, we have that

$$\langle \tilde{d}_i(v), \tilde{d}_j(v) \rangle = \|\tilde{d}_i(v)\|^2 \delta_{ij} \quad \text{and} \quad \langle \tilde{s}_i(g), \tilde{s}_j(g) \rangle = \|\tilde{s}_i(g)\|^2 \delta_{ij}$$
(18)

where δ_{ij} is the Kronecker delta function, i.e., $\delta_{ij} = 1$, if i = j; $\delta_{ij} = 0$, otherwise.

¹¹The usual SVD yields unit-norm vectors $\tilde{d}_i(v)$, $\tilde{s}_i(g)$ and additional singular values σ_i . In this paper, what we are really interested is the fixed rank approximation of $\hat{R}(P)$ via SVD. Therefore, once SVD is computed, one can lump σ_i into either $\tilde{d}_i(v)$ or $\tilde{s}_i(g)$ or even divide $\tilde{d}_i(v)$ and multiply $\tilde{s}_i(g)$ by some constant simultaneously without changing the decomposition, since $\tilde{d}_i(v)$ and $\tilde{s}_i(g)$ appear together in a product in the decomposition.

3 Estimation of shape and radiance for non-Lambertian scenes

In this section we present our algorithm to recover the representation of shape and radiance described in the previous section from a collection of images.

3.1 Shape estimation

Shape, in our context, is described by a representation of the surface S relative to any Euclidean reference frame. When S is represented explicitly, one can look for the solution \hat{S} via a local descent along the gradient of E. The optimality condition associated with equation (11) is given in the following theorem.

Theorem 1 (optimality condition). Let $\Phi_{\mathbf{X}}$, Φ_N be the first-order derivatives of Φ with respect to \mathbf{X} and N and $\Phi_{\mathbf{X}N}$, Φ_{NN} be the second-order derivatives. We assume that Φ_{NN} can be decomposed as: $\Phi_{NN} = \sum_{i=1}^{k} \lambda_i p_i p_i^T$ where $\lambda_i \in \mathbb{R}$ and $p_i \in \mathbb{R}^3$ (note that this decomposition is always possible since Φ_{NN} is real and symmetric). We have that the following partial differential equation is the gradient descent flow for the cost (11):

$$S_t = \left(2H\Phi - \langle \Phi_{\mathbf{X}}, N \rangle - 2H \langle \Phi_N, N \rangle - trace(\Phi_{\mathbf{X}N}) + N^T \Phi_{\mathbf{X}N} N + \sum_{i=1}^k \lambda_i \mathbf{II}(P_N^{\perp} p_i)\right) N.$$
(19)

where P_N^{\perp} is the projection from \mathbb{R}^3 to $T_P(S)$, i.e., $P_N^{\perp} = I - NN^T$, H is the mean curvature and $\mathbf{II}(t)$ is the second fundamental form of a vector $t \in T_P(S)$.

Proof. See Appendix A.

Note that equation (19) involves second-order derivatives: $\Phi_{\mathbf{X}N}$ and Φ_{NN} and no higher-order derivatives. This should not be surprising because the cost functional involves N in the integrand, which is the first-order variation of the surface S. In practice the following flow based on the first-order derivatives in equation (19) yields similar results to that of the flow (19), while saving a significant amount of computations.

$$S_t = \left(2H\Phi - \langle \Phi_{\mathbf{X}}, N \rangle - 2H \langle \Phi_N, N \rangle \right) N.$$
(20)

The calculation of the flow above reveals some interesting structure, as major simplification occur after equations (17) and (18).

We will prove a stronger result than needed to compute flow (20). In particular, we will show that even if the dimension of \mathbb{A}_m^r is higher than 1, i.e., $r \ge 1$, the resulting flow still takes a simple expression. Let r be the rank the residual matrix defined in equation (13). Suppose that the columns of R(P) satisfy an affine subspace constraint of dimension r_0 and we take r_0 terms from the SVD of $\widehat{R}(P)$ (equation (14)). Therefore, the function Φ takes the expression

$$\Phi(P) = \|\tilde{d}_{r_0+1}(v)\tilde{s}_{r_0+1}^T(g) + \tilde{d}_{r_0+2}(v)\tilde{s}_{r_0+2}^T(g) + \dots + \tilde{d}_r(v)\tilde{s}_r^T(g)\|_F^2.$$
(21)

Let ϕ_{ij} be the (i, j)-th element of $\Phi(P)$.

Theorem 2 (differentiation of the score). Let ξ indicate the arguments of Φ , i.e., ξ is one of X_1, X_2, X_3 , N_1, N_2, N_3 . Then

$$\dot{\Phi} = \sum_{i,j=1}^{m,n} 2\phi_{ij}\dot{\tilde{R}}_{ij}$$
(22)

where the dot indicates differentiation with respect to ξ .

Proof. We will first show that $s_i(g)$ sums to zero for any $1 \le i \le r$. Recall that the columns of the

residual matrix $\widehat{R}(P)$ sum to zero (equation (17)). We have that

$$\widehat{R}(P)\mathbf{1}_n = \sum_{k=1}^r \widetilde{d}_k(v) \left\langle \widetilde{s}_k(g), \mathbf{1}_n \right\rangle = 0$$
(23)

Multiplying both sides of equation (23) on the left by $\tilde{d}_i^T(v)$ where $1 \le i \le r$, we have

$$\tilde{d}_i^T(v)\sum_{k=1}^r \tilde{d}_k(v)\left<\tilde{s}_k(g), \mathbf{1}_n\right> = \|\tilde{d}_i(v)\|^2\left<\tilde{s}_i(g), \mathbf{1}_n\right> = 0.$$
(24)

Since $\|\tilde{d}_i(v)\|^2 \neq 0$, we have

$$\langle \tilde{s}_i(g), \mathbf{1}_n \rangle = 0, \quad 1 \le i \le r.$$
(25)

We define

$$\phi^{i} \doteq \tilde{R}^{i} - \tilde{d}_{0}(v_{i})\mathbf{1}_{n} - \sum_{k=1}^{r_{0}} \tilde{d}_{k}(v_{i})\tilde{s}_{k}(g), \quad i = 1, 2, \dots, m,$$

$$\phi_{j} \doteq \tilde{R}_{j} - \tilde{d}_{0}(v) - \sum_{k=1}^{r_{0}} \tilde{d}_{k}(v)\tilde{s}_{k}(g_{j}), \quad j = 1, 2, \dots, n,$$

where \tilde{R}^i is the *i*-th row of \tilde{R} and \tilde{R}_j is the *j*-th column of \tilde{R} , i.e., $\tilde{R}^i = [\tilde{R}_{i1}, \tilde{R}_{i2}, \dots, \tilde{R}_{in}]^T$ and $\tilde{R}_j = [\tilde{R}_{1j}, \tilde{R}_{2j}, \dots, \tilde{R}_{mj}]^T$. Expanding the derivative we get

$$\dot{\Phi} = \sum_{i,j=1}^{n,m} \dot{\phi}_{ij}^2 = \sum_{i,j=1}^{n,m} 2\phi_{ij} \left(\dot{\tilde{R}}_{ij} - \dot{\tilde{d}}_0(v_i) - \sum_{k=1}^{r_0} \dot{\tilde{d}}_k(v_i) \tilde{s}_k(g_j) - \sum_{k=1}^{r_0} \tilde{d}_k(v_i) \dot{\tilde{s}}_k(g_j) \right) \\
= \sum_{i,j=1}^{n,m} 2\phi_{ij} \dot{\tilde{R}}_{ij} - \sum_{i=1}^{n} \dot{\tilde{d}}_0(v_i) \left\langle \phi^i, \mathbf{1}_n \right\rangle - \sum_{k=1}^{r_0} \sum_{i=1}^{n} \dot{\tilde{d}}_k(v_i) \left\langle \phi^i, \tilde{s}_k(g) \right\rangle - \sum_{k=1}^{r_0} \sum_{j=1}^{m} \dot{\tilde{s}}_k(g_j) \left\langle \phi_j, \tilde{d}_k(v) \right\rangle.$$

However, from equations (26) and (26) we see that ϕ^i is in the span of $\tilde{s}_{r_0+1}(g), \tilde{s}_{r_0+2}(g), \ldots, \tilde{s}_r(g)$ and ϕ_j is in the span of $\tilde{d}_{r_0+1}(v), \tilde{d}_{r_0+2}(v), \ldots, \tilde{d}_r(v)$. Therefore, from equation (25) we have $\langle \phi^i, \mathbf{1}_n \rangle = 0$, and together with equations (18), we can conclude the proof by noting that the only term that contributes to the derivative is $\sum_{i,j=1}^{n,m} 2\phi_{ij}\dot{\tilde{R}}_{ij}$.

As a consequence of the previous result, flow (20) for arbitrary dimensional subspace in an explicit form read as:

$$S_t = \left(2H\Phi - \sum_{i,j=1}^{m,n} 2\phi_{ij} \left\langle \frac{\partial \tilde{R}_{ij}}{\partial \mathbf{X}} + 2H \frac{\partial \tilde{R}_{ij}}{\partial N}, N \right\rangle \right) N.$$
(26)

3.1.1 Implementation notes

To compute ϕ_{ij} we need to obtain the first singular value and the corresponding singular vectors of the residual matrix $\widehat{R}(P)$ (or, equivalently, the rank-1 approximation of $\widehat{R}(P)$). Since we are only interested in the rank-1 approximation, we do not have to carry a full singular value decomposition. Instead, we use the power iteration [9] to efficiently compute the largest singular value and its corresponding singular vectors. Note that the algorithm presented in [15] uses the rank-2 constraint and therefore requires to compute the first two singular values and the corresponding singular vectors. In general the algorithm presented in this paper is about twice faster than the one in [15] because the computation of singular values and vectors at every point dominates both algorithms.

We implement the flow (26) using level set methods [27]. Naturally, as with most of these variational techniques, one can only hope to achieve convergence to a local extremum of the original cost functional,

since the flow is based on the gradient descent principle, and existence and uniqueness results are not available for this class of flows. In the experimental section we will give empirical validation to this approach by testing the flow above on real image data starting from generic initial conditions.

3.2 Radiance estimation

Once the surface \hat{S} has been found, one can use the representation of the radiance to generate images by "radiance-mapping" the tensor R(P) onto the surface S. Naturally, the visualization of S in this case is view-dependent, since different columns of R(P) contribute to the image of the same point P depending on the viewpoint g_i .

The radiance map is provided by the functions $d_0(v)$, $d_1(v)$ and $s_1(g)$, estimated at each point of the surface, P, using the singular value decomposition of the measured radiance tensor \tilde{R} , according to Corollary 1 and equation (15). Given a novel vantage point g', the corresponding function $s_1(g')$ can be interpolated from the existing $s_1(g_j)$. One simple way of doing so is to find the three views closest to g', and then use the linear interpolation to obtain $s_1(g')$ from $s_1(g_j)$. This technique also allows extrapolating the radiance; as we show in the experimental section, one can notice artifacts when comparing the results to actual images obtained from a novel viewpoint. However, such artifacts are only noticeable by direct comparison. Notice that $d_0(v)$ and $d_1(v)$ do not depend on the viewpoint, and therefore do not need to be interpolated. Since $s_1(g')$ is linearly interpolated from $s_1(g_j)$, this new radiance component does not increase the dimension of the subspace spanned by the columns of the radiance tensor and therefore is consistent with the affine subspace constraint (Proposition 1).

Notice that the images generated from the radiance map are significantly different than those generated by "texture mapping" the images \tilde{I} onto the surface S. In fact, the function $s_1(g)$ depends directly on the viewpoint, and therefore when the viewpoint moves, the highlights move on the estimated surface, giving an overall result that is visually comparable with image-based rendering techniques that assume true surface shapes [10, 5].

4 Experiments

In this section we report the experimental results of our algorithm tested on three datasets: "Van Gogh", "Buddha" and "elephant". The first two datasets (shown in Figure 1) are courtesy of Jean-Yves Bouguet and Radek Grzeszczuk (Intel Corp.). The third dataset (shown in the top row of Figure 10) is courtesy of Daniel Wood (University of Washington). The Van Gogh statue is made of polished metal, and is highly specular. There is a total of 339 images in the dataset. Pseudo-ground truth has been generated by laser or shadow scanning followed by mesh polishing (Figure 4). Buddha is actually a synthetic scene. There is a total of 281 images in the dataset. Ground truth is available (Figure 7). In Figure 4 we show the estimates of shape produced by the algorithm described in Section 3.1, together with the estimates obtained with the algorithm of [16], both compared with pseudo ground truth. In both algorithms, the numerical grids we use are of size $128 \times 128 \times 128$. Our estimate is obviously not as crisp as the ground truth, but it does capture important details on the face. Figure 8 shows the evolution of the estimate of shape. In Figure 9 we show synthetic images generated using the radiance map, as described in Section 3.2. Note that the specularities move with the viewpoint. In Figure 5 we show a few synthetic images compared with the real images from the same vantage point. In Figure 7 we show the estimated shape for the Buddha in Figure 1. The numerical grid size is $128 \times 128 \times 128$. In this case, ground truth is available since the images are synthetic. We also show the results obtained with the algorithm of [16]. In Figure 6 we show images synthesized from the model, compared with corresponding true images. In



Figure 4: Estimated shape (top), compared with pseudo-ground truth (bottom), obtained with a 3D laser scanner and mesh polishing. Our results improve those obtained with the algorithm of [16] (middle).

Figure 8 we show the evolution of shape, and in Figure 9 we show several novel views. In Figure 10 (top row) we show several views of an elephant made of polished marble. There is a total of 397 images in the dataset. The numerical grid size we used is again $128 \times 128 \times 128$. The estimated shape of our algorithms is reported in Figure 11 compared with pseudo-ground truth and that obtained using [16]. In the bottom row of Figure 10 we show images synthesized from the model, whose viewing positions and directions are the same as those in the top row.

In Table 1 we summarize the shape error for different approximations of surface reflectance. The error is measured by the ratio between the volume of the symmetric difference between the estimated shape and the true shape (or the pseudo-ground truth) and the volume of the true shape (or the pseudo-ground



Figure 5: (COLOR) Synthetic images using the estimated radiance tensor (top) compared with the true images taken from the same vantage point. Note that one can actually read the text at the base of the bust. This is obtained from the radiance estimate, not from texture mapping.

truth). Let ψ be the level set function for surface S. Suppose ψ is negative inside S and positive outside

S. Then the volume contained by S, defined as S_{in} , can be measured as

$$\operatorname{Vol}(S_{in}) = \int_{\mathbb{R}^3} (1 - \mathcal{H}(\psi)) dx dy dz$$
(27)

where $\mathcal{H}(x) : \mathbb{R} \to \{0,1\}$ is the Heaviside function: $\mathcal{H}(x) = 1$ if $x \ge 0$; $\mathcal{H}(x) = 0$ otherwise. In practice, one can mollify the Heaviside function with a smooth approximation [4]. The volume of the

Reflectance models	Van Gogh	Buddha	elephant
Lambertian (the algorithm presented in [16])	6.9%	5.5%	24.3%
1-dimensional affine subspace (the proposed al-	5.7%	3.5%	7.3%
gorithm)			
2-dimensional affine subspace (a modified ver-	5.6%	3.4%	7.2%
sion of the proposed algorithm by using the first			
and second singular values and vectors in ap-			
proximating $\widehat{R}(P)$ in equation (14))			

Table 1: Shape error comparison chart for different reflectance approximations. The error is measured by the ratio between the volume of the symmetric difference between the estimated shape and the true shape (or the pseudo-ground truth) and the volume of the true shape (or the pseudo-ground truth). We observe that using the proposed one-dimensional affine subspace constraint for the radiance tensor, we can reduce the shape error by a factor of 2 in average, while using higher-order affine subspaces does not improve the results much.

symmetric difference between S_{in} and T_{in} can be calculated by:

$$\operatorname{Vol}(S_{in}\Delta T_{in}) = \operatorname{Vol}(S_{in}) + \operatorname{Vol}(T_{in}) - 2\operatorname{Vol}(S_{in} \cap T_{in}).$$
(28)

We observe that using the proposed one-dimensional affine subspace constraint for the radiance tensor, we can reduce the shape error by a factor of 2 in average, while using higher-order affine subspaces does not improve the results much. In Table 2 we show the degradation of the reconstruction as a function of the size of the patch Ω_P for the Vangogh dataset. We tested neighborhood sizes from 3×3 to 19×19 . We use odd sizes to have the neighborhoods symmetric around the center point. The unit of the neighborhood size is chosen to be corresponding to the actual pixel size in the best view, for instance 5×5 means that the projected neighborhood in the best view occupies an approximate 5×5 region in image pixels. We observe that the proposed algorithm is very robust with respect to the neighborhoods size in the sense that the reconstruction errors are almost the same from 7×7 to 15×15 neighborhoods. When the neighborhood is chosen too small, the algorithm is sensitive to image noise and therefore has

Size of Ω_P	3×3	5×5	7×7	9×9	11×11	13×13	15×15	17×17	19×19
Shape error	14.8%	7.6%	6.3%	6.0%	5.7%	5.8%	6.1%	6.4%	6.6%

Table 2: Shape error comparison chart for different sizes of Ω_P for the Vangogh dataset. The error is measured by the ratio between the volume of the symmetric difference between the estimated shape and the pseudo-ground truth and the volume of the pseudo-ground truth. The unit of the neighborhood size is chosen to be corresponding to the actual pixel size in the best view, for instance 5×5 means that the projected neighborhood in the best view will occupy an approximate 5×5 region in image pixels.

trouble to converge. When the neighborhood is chosen too big, the algorithm has trouble to capture sharp features presented in the object shape.

Occlusions are handled by computing visibility at each step of the iteration. Therefore, the technique we present is computationally intensive and processing an entire dataset takes several hours. On the other hand, the algorithm requires no manual intervention, no intermediate step, no mesh polishing and no texture mapping after reconstruction. Therefore, its computational cost should be compared to implementing the entire pipeline from images to rendering.



Figure 6: (COLOR) Synthetic images obtained with the estimated radiance tensor field (top) compared with the true images taken from the same vantage point.



Figure 7: Estimated shape (top), compared with ground truth (bottom), also compared with the results obtained by the algorithm of [16] (middle).

5 Discussion

We have presented a novel algorithm for estimating dense shape and non-Lambertian photometry from a collection of images. Our algorithm relies on an affine subspace constraint on the radiance tensor field, which is derived from the diffuse+specular reflection model commonly used in Computer Graphics, in the sense elucidated in Proposition 1. While one could dismiss the analysis and just introduce the cost function (9) point-blank without detracting from the algorithm proposed (which is validated experimentally), the proposition indicates precisely under what conditions the affine subspace constraint is satisfied, i.e., what the underlying *mathematical model* is. Naturally, the closest the scene is to satisfying



Figure 8: Shape evolution for Van Gogh (top) and Buddha (bottom).

the assumptions (i.e. the closest it is to smooth shape, diffuse+specular reflection, fixed distant illumination) the smaller dimension the subspace spanned by the columns of $\tilde{R}(P)$ is. However, even though only ideal scenes viewed from noiseless images will satisfy the assumptions exactly, we can still exploit the discrepancy derived from the idealized model to define a constraint that can be used to reconstruct the scene from real images.

Those that object to the restrictiveness of the model laid out in Proposition 1 will be relieved to know that extension to higher dimensional affine subspace is conceptually and computationally trivial. One will need to take more terms from the SVD, but Theorem 1 assures that the gradient flow can be computed essentially in the same way. However, it can be verified experimentally that, for most scenes, an increase in the dimension of the model does not yield a significant improvement in the reconstruction, further validating the mathematical model proposed (see Table 1).

Our algorithm can handle sharp changes of the radiance profile: In Figure 5, one can actually *read* the text at the base of the bust from the reconstructed radiance. Note that there is no restriction whatsoever imposed on the variation of the diffuse and specular components of the radiance, and nowhere it is



Figure 9: (COLOR) Synthetic images obtained from the estimated radiance. As it can be seen, the appearance changes significantly with the vantage point.

assumed that it be constant or smooth. What is assumed to be constant is the surface roughness, *not the albedo*, so we can handle heavily textured objects. On the other hand, our algorithm does not *require* strong texture or point features to be visible, and returns a dense estimate of shape, with no need to interpolate or triangulate a surface from sparse points.

Note also that, although the measured radiance tensor at a given point P is assembled using a local approximation of the surface with the tangent plane T_PS , this does not mean that our algorithm only works for planar surfaces: In fact, the radiance tensor at a nearby point Q is computed using the tangent plane T_QS that is not constrained to be similar to T_PS . If one thinks of R(P) as a "signature" attached to $P \in S$, the model imposes no constraint that nearby points should have similar signatures.

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Figure 10: Top row: three images from the elephant dataset (courtesy of Daniel Wood, University of Washington). Bottom row: synthetic views generated using the estimated radiance. The structure and position of specular highlights is correctly captured; there are some visualization artifacts at the boundaries, but note that even the text on the small label is visible on the left image. Note that this is an estimate of the radiance, not a texture map.



Figure 11: Estimated shape (top row) of the scene in Figure 10, compared with pseudo-ground truth (bottom row), obtained with a 3D laser scanner and mesh polishing. Our results improve those obtained with [16] (middle row). The ear is not clear in the reconstruction, although it is well captured as radiance (Figure 10).

A Gradient descent flow for $\int_{S} \Phi(\mathbf{X}, N) dA$

In this section we present the proof for Theorem 1, i.e. deriving the gradient descent flow for the cost (11). Note that the analysis for the cost (11) was also done by Faugeras and Keriven in [8] and can be found in [19] (in French). In particular, Faugeras and Keriven derived the Euler-Lagrange equations for the cost functional and then designed a flow based on it to find the optimal shape. However, in their derivation it is not immediate to see that the resulting flow minimizes the cost functional. In this paper we present a way of directly minimizing the cost functional (11) and show that the flow considered by Faugeras and Keriven in [8] is indeed the gradient descent for the cost (11).

We shall express the surface S using a fixed parameterization $\mathbf{X} : U \to S$; $(u, v) \mapsto \mathbf{X}(u, v)$. We will further assume that U is simply connected to exploit integration by parts. This seems like a restrictive assumption to start with but, in the end, the resulting flow will have a completely local expression which does not depend on the choice of the parameterization. Let E, F, and G be the coefficients of the first fundamental form of S under the parameterization \mathbf{X} and e, f, and g be the coefficients of the second fundamental form, i.e., $E \doteq \langle \mathbf{X}_u, \mathbf{X}_u \rangle$, $F \doteq \langle \mathbf{X}_u, \mathbf{X}_v \rangle$, $G \doteq \langle \mathbf{X}_v, \mathbf{X}_v \rangle$, and $e \doteq \langle N, \mathbf{X}_{uu} \rangle$, $f \doteq$ $\langle N, \mathbf{X}_{uv} \rangle$, $g \doteq \langle N, \mathbf{X}_{vv} \rangle$. We can write the area form dA using the coefficients of the first fundamental form $dA = \sqrt{EG - F^2} du dv$. By definition $N = \frac{\mathbf{X}_u \times \mathbf{X}_v}{\|\mathbf{X}_u \times \mathbf{X}_v\|} = \frac{\mathbf{X}_u \times \mathbf{X}_v}{\sqrt{EG - F^2}}$. We will also use the following lemma from linear algebra, which can be verified by direct substitution:

Lemma 1. Let $\mathbf{x}, \mathbf{y} \in \mathbb{R}^3$ and $A \in \mathbb{R}^{3 \times 3}$. We have the following identity:

$$\mathbf{x} \times (A\mathbf{y}) + (A\mathbf{x}) \times \mathbf{y} = trace(A)(\mathbf{x} \times \mathbf{y}) - A^{T}(\mathbf{x} \times \mathbf{y}).$$
(29)

The general scheme we choose to minimize cost (11) is to first augment the parameterization with an additional time variable t: $\mathbf{X}(u, v, t)$ and then update \mathbf{X} (and thus evolve S) according to a partial differential equation $S_t = \mathcal{F}N$ starting from a generic shape (such as a large sphere). We shall show that the choice of the partial differential equation made in Theorem 1 minimizes cost (11). We use \mathbf{X}_t to indicate the derivative of \mathbf{X} with respect to time and replace \mathbf{X}_t with S_t when the whole expression does not depend on the parameterization. Throughout our computation, we will always omit purely tangential terms whenever they appear inside an inner product with the surface variation \mathbf{X}_t or S_t , because they will not change the shape of S [6]. We will first prove the following theorem, which will be used in the further computation and enlighten it.

Theorem 3 (mean curvature flow). Let S be a closed regular surface in \mathbb{R}^3 . The gradient flow to minimize the surface area $\int_S dA$ is the mean curvature flow:

$$S_t = HN. ag{30}$$

Proof. Computing the time derivative of the surface area, we obtain

$$\begin{split} \frac{\partial}{\partial t} \int_{S} dA &= \iint \sqrt{EG - F^{2}}_{t} du dv = \iint \frac{E_{t}G + EG_{t} - 2FF_{t}}{2\sqrt{EG - F^{2}}} du dv \\ &= \iint \frac{G \left\langle \mathbf{X}_{ut}, \mathbf{X}_{u} \right\rangle + E \left\langle \mathbf{X}_{vt}, \mathbf{X}_{v} \right\rangle - F\left(\left\langle \mathbf{X}_{vt}, \mathbf{X}_{u} \right\rangle + \left\langle \mathbf{X}_{ut}, \mathbf{X}_{v} \right\rangle\right)}{\sqrt{EG - F^{2}}} du dv \\ &= \iint -\frac{\left\langle G\mathbf{X}_{uu}, \mathbf{X}_{t} \right\rangle + E \left\langle \mathbf{X}_{vv}, \mathbf{X}_{t} \right\rangle - 2F \left\langle \mathbf{X}_{uv}, \mathbf{X}_{t} \right\rangle}{\sqrt{EG - F^{2}}} du dv \\ &= \iint -\frac{\left\langle G \left\langle \mathbf{X}_{uu}, N \right\rangle N + E \left\langle \mathbf{X}_{vv}, N \right\rangle N - 2F \left\langle \mathbf{X}_{uv}, N \right\rangle N, \mathbf{X}_{t} \right\rangle}{\sqrt{EG - F^{2}}} du dv \\ &= \iint \left\langle -\frac{Ge + Eg - 2Ff}{EG - F^{2}} N, \mathbf{X}_{t} \right\rangle \sqrt{EG - F^{2}} du dv \\ &= \iint \left\langle -2HN, S_{t} \right\rangle dA. \end{split}$$

Note that for the sake of simplicity, we have omitted the integration domain for u and v. If we choose $S_t = HN$, then we have

$$\frac{\partial}{\partial t}\int_{S}dA=\int\int\!\!\int-2H^{2}dA\leq0$$

Therefore, $S_t = HN$ minimizes the surface area.

Now we come to our original cost functional $\int_{S} \Phi(\mathbf{X},N) dA.$

Proof of Theorem 1.

$$\frac{\partial}{\partial t} \int_{S} \Phi(\mathbf{X}, N) dA = \iint \frac{\partial}{\partial t} \left(\Phi(\mathbf{X}, N) \sqrt{EG - F^{2}} \right) du dv$$

$$= \iint \left(\langle \Phi_{\mathbf{X}}, \mathbf{X}_{t} \rangle + \langle \Phi_{N}, N_{t} \rangle \right) \sqrt{EG - F^{2}} + \Phi \sqrt{EG - F^{2}}_{t} du dv$$

$$= \iint \left(\left(\langle \Phi_{\mathbf{X}}, N \rangle - 2H\Phi \right) \langle N, \mathbf{X}_{t} \rangle + \langle \Phi_{N}, N_{t} \rangle \right) \sqrt{EG - F^{2}} du dv$$

$$= \iint \left(\langle \Phi_{\mathbf{X}}, N \rangle - 2H\Phi \right) \langle N, S_{t} \rangle dA$$

$$+ \iint \left\langle \Phi_{N}, \mathbf{X}_{ut} \times \mathbf{X}_{v} + \mathbf{X}_{u} \times \mathbf{X}_{vt} - \frac{\mathbf{X}_{u} \times \mathbf{X}_{v}}{\sqrt{EG - F^{2}}} \sqrt{EG - F^{2}}_{t} \right\rangle du dv$$

$$= \iint \left(\langle \Phi_{\mathbf{X}}, N \rangle - 2H\Phi \right) \langle N, S_{t} \rangle dA$$

$$+ \iint \left\langle \Phi_{N}, \mathbf{X}_{ut} \times \mathbf{X}_{v} + \mathbf{X}_{u} \times \mathbf{X}_{vt} \right\rangle \langle N, S_{t} \rangle dA$$

$$+ \iint \left\langle \Phi_{N}, \mathbf{X}_{ut} \times \mathbf{X}_{v} + \mathbf{X}_{u} \times \mathbf{X}_{vt} \right\rangle \langle N, S_{t} \rangle dA$$

$$+ \iint \left\langle \Phi_{N}, \mathbf{X}_{ut} \times \mathbf{X}_{v} + \mathbf{X}_{u} \times \mathbf{X}_{vt} \right\rangle du dv$$
(31)

We continue on the second term $\iint \langle \Phi_N, \mathbf{X}_{ut} \times \mathbf{X}_v + \mathbf{X}_u \times \mathbf{X}_{vt} \rangle dudv$:

$$\begin{split} &\iint \langle \Phi_N, \mathbf{X}_{ut} \times \mathbf{X}_v + \mathbf{X}_u \times \mathbf{X}_{vt} \rangle \, du dv = \iint \langle \mathbf{X}_v \times \Phi_N, \mathbf{X}_{ut} \rangle + \langle \Phi_N \times \mathbf{X}_u, \mathbf{X}_{vt} \rangle \, du dv \\ &= \iint - \langle \mathbf{X}_{uv} \times \Phi_N + \mathbf{X}_v \times (\Phi_N)_u + (\Phi_N)_v \times \mathbf{X}_u + \Phi_N \times \mathbf{X}_{uv}, \mathbf{X}_t \rangle \, du dv \\ &= \iint - \langle \mathbf{X}_v \times (\Phi_N)_u + (\Phi_N)_v \times \mathbf{X}_u, \mathbf{X}_t \rangle \, du dv \\ &= \iint - \langle \mathbf{X}_v \times (\Phi_{N\mathbf{X}}\mathbf{X}_u + \Phi_{NN}N_u) + (\Phi_{N\mathbf{X}}\mathbf{X}_v + \Phi_{NN}N_v) \times \mathbf{X}_u, \mathbf{X}_t \rangle \, du dv \\ &= \iint - \langle \mathbf{X}_v \times (\Phi_{N\mathbf{X}}\mathbf{X}_u) + (\Phi_{N\mathbf{X}}\mathbf{X}_v) \times \mathbf{X}_u, \mathbf{X}_t \rangle \, du dv \\ &= \iint \langle \mathbf{X}_v \times (\Phi_{N\mathbf{X}}\mathbf{X}_u) + (\Phi_{N\mathbf{X}}\mathbf{X}_v) \times \mathbf{X}_u, \mathbf{X}_t \rangle \, du dv \end{split}$$

We can simplify the first term as follows:

$$\iint - \langle \mathbf{X}_{v} \times (\Phi_{N\mathbf{X}}\mathbf{X}_{u}) + (\Phi_{N\mathbf{X}}\mathbf{X}_{v}) \times \mathbf{X}_{u}, \mathbf{X}_{t} \rangle \, dudv$$

$$= \iint \langle \operatorname{trace}(\Phi_{N\mathbf{X}})(\mathbf{X}_{u} \times \mathbf{X}_{v}) - \Phi_{N\mathbf{X}}^{T}(\mathbf{X}_{u} \times \mathbf{X}_{v}), \mathbf{X}_{t} \rangle \, dudv$$

$$= \iint (\operatorname{trace}(\Phi_{\mathbf{X}N}) - N^{T}\Phi_{\mathbf{X}N}N) \, \langle N, S_{t} \rangle \, dA. \tag{32}$$

We can also simplify the second term by noting that Φ_{NN} can be decomposed as follows:

$$\Phi_{NN} = \sum_{i=1}^{k} \lambda_i p_i p_i^T \tag{33}$$

where $\lambda_i \in \mathbb{R}$ and $p_i \in \mathbb{R}^3$. We wish to make two remarks. First, as a real 3×3 symmetric matrix, Φ_{NN} can always be decomposed into the form of equation (33) (for instance via the *eigenvalue decomposition*). Second, equation (33) is more general than the eigenvalue decomposition, because it allows k to

take any number and p_i is not necessarily orthogonal to each other. Thus, we have:

$$\begin{split} &\iint - \langle \mathbf{X}_{v} \times (\Phi_{NN}N_{u}) + (\Phi_{NN}N_{v}) \times \mathbf{X}_{u}, \mathbf{X}_{t} \rangle \, du dv \\ &= \iint - \sum_{i=1}^{k} \lambda_{i} \left\langle \left\langle p_{i}, N_{u} \right\rangle \mathbf{X}_{v} \times p_{i} + \left\langle p_{i}, N_{v} \right\rangle p_{i} \times \mathbf{X}_{u}, \mathbf{X}_{t} \right\rangle \, du dv \\ &= \iint - \sum_{i=1}^{k} \lambda_{i} \left\langle p_{i}, N_{u} \right\rangle \left\langle \mathbf{X}_{v} \times p_{i}, N \right\rangle + \left\langle p_{i}, N_{v} \right\rangle \left\langle p_{i} \times \mathbf{X}_{u}, N \right\rangle \left\langle N, \mathbf{X}_{t} \right\rangle \, du dv \\ &= \iint - \sum_{i=1}^{k} \lambda_{i} \left\langle P_{N}^{\perp} p_{i}, [N_{u}, N_{v}] \begin{bmatrix} \left\langle P_{N}^{\perp} p_{i}, N \times \mathbf{X}_{v} \right\rangle \\ \left\langle P_{N}^{\perp} p_{i}, \mathbf{X}_{u} \times N \right\rangle \end{bmatrix} \right\rangle \, du dv \end{split}$$

The last step follows from the fact that N_u , N_v , $\mathbf{X}_u \times N$ and $N \times \mathbf{X}_v$ are all tangent vectors. If we construct a curve $C(t) : (-\epsilon, \epsilon) \to S$ with $C_t = P_N^{\perp} p_i$, we can easily verify that

$$u_t = \sqrt{EG - F^2} \left\langle P_N^{\perp} p_i, N \times \mathbf{X}_v \right\rangle$$
 and $v_t = \sqrt{EG - F^2} \left\langle P_N^{\perp} p_i, \mathbf{X}_u \times N \right\rangle$,

and

$$\left\langle P_{N}^{\perp}p_{i}, [N_{u}, N_{v}] \begin{bmatrix} \left\langle P_{N}^{\perp}p_{i}, N \times \mathbf{X}_{v} \right\rangle \\ \left\langle P_{N}^{\perp}p_{i}, \mathbf{X}_{u} \times N \right\rangle \end{bmatrix} \right\rangle = \sqrt{EG - F^{2}} \left\langle P_{N}^{\perp}p_{i}, N_{u}u_{t} + N_{v}v_{t} \right\rangle = \sqrt{EG - F^{2}} \mathbf{II}(P_{N}^{\perp}p_{i}).$$

Hence we have

$$\iint -\langle \mathbf{X}_v \times (\Phi_{NN}N_u) + (\Phi_{NN}N_v) \times \mathbf{X}_u, \mathbf{X}_t \rangle \, du dv = \iint -\sum_{i=1}^k \lambda_i \mathbf{II}(P_N^{\perp}p_i) dA.$$
(34)

Putting equations (31), (32) and (34) together, we arrive at

$$\frac{\partial}{\partial t} \int_{S} \Phi(\mathbf{X}, N) dA = \int_{S} \left(\langle \Phi_{\mathbf{X}}, N \rangle - 2H\Phi + 2H \langle \Phi_{N}, N \rangle + (\operatorname{trace}(\Phi_{\mathbf{X}N}) - N^{T}\Phi_{\mathbf{X}N}N) - \sum_{i=1}^{k} \lambda_{i} \mathbf{II}(P_{N}^{\perp}p_{i}) \right) \langle N, S_{t} \rangle \, dA.$$
(35)

It is immediate that the gradient descent flow is given by

$$S_t = \left(2H\Phi - \langle \Phi_{\mathbf{X}}, N \rangle - 2H \langle \Phi_N, N \rangle - \operatorname{trace}(\Phi_{\mathbf{X}N}) + N^T \Phi_{\mathbf{X}N} N + \sum_{i=1}^k \lambda_i \mathbf{II}(P_N^{\perp} p_i)\right) N.$$
(36)

Remark 4. Even though in the computation we have ignored all tangential terms that appear inside an inner product with \mathbf{X}_t or S_t , it is possible, although far more laborious, to show that they indeed cancel out through a more careful derivation.

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