

Parallelizable global conformal parameterization of simply-connected surfaces via partial welding*

Gary P. T. Choi[†], Yusan Leung-Liu[‡], Xianfeng Gu[§], and Lok Ming Lui[¶]

Abstract.

Conformal surface parameterization is useful in graphics, imaging and visualization, with applications to texture mapping, atlas construction, registration, remeshing and so on. With the increasing capability in scanning and storing data, dense 3D surface meshes are common nowadays. While meshes with higher resolution better resemble smooth surfaces, they pose computational difficulties for the existing parameterization algorithms. In this work, we propose a novel parallelizable algorithm for computing the global conformal parameterization of simply-connected surfaces via partial welding maps. A given simply-connected surface is first partitioned into smaller subdomains. The local conformal parameterizations of all subdomains are then computed in parallel. The boundaries of the parameterized subdomains are subsequently integrated consistently using a novel technique called partial welding, which is developed based on conformal welding theory. Finally, by solving the Laplace equation for each subdomain using the updated boundary conditions, we obtain a global conformal parameterization of the given surface, with bijectivity guaranteed by quasi-conformal theory. By including additional shape constraints, our method can be easily extended to achieve disk conformal parameterization for simply-connected open surfaces and spherical conformal parameterization for genus-0 closed surfaces. Experimental results are presented to demonstrate the effectiveness of our proposed algorithm. When compared to the state-of-the-art conformal parameterization methods, our method achieves a significant improvement in both computational time and accuracy.

Key words. Conformal parameterization, conformal welding, parallelization, simply-connected surface

AMS subject classifications. 65D18, 68U05, 68U10, 68W10

1. Introduction. From mobile games to high-resolution movies, from small 3D printed desk toys to large aircraft engines, 3D geometric models are everywhere nowadays. With the advancement of computing technologies and scanning devices, 3D geometric data can be created or acquired easily. However, at the same time, the scale of the data grows rapidly. In many situations, it is necessary to handle dense geometric data with hundreds of thousands, or even millions of vertices.

In geometry processing, a common representation of 3D objects is triangulated 3D surfaces. To simplify various tasks that are to be performed on the 3D surfaces, one possible way is to transform the 3D surfaces into a simpler 3D shape or a 2D shape. This process is known as surface parameterization. With the aid of surface parameterization, we can perform the tasks on the simpler domain and transform the results back to the original 3D surfaces instead of working on them directly. For instance, under surface parameterization, PDEs on complicated

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[†]John A. Paulson School of Engineering and Applied Sciences, Harvard University, Cambridge, MA 02138, USA (pchoi@g.harvard.edu).

[‡]Department of Mathematics, The Chinese University of Hong Kong, Hong Kong (ylleung@math.cuhk.edu.hk).

[§]Department of Computer Science, Stony Brook University, Stony Brook, NY 11794, USA (gu@cs.stonybrook.edu).

[¶]Department of Mathematics, The Chinese University of Hong Kong, Hong Kong (lmhui@math.cuhk.edu.hk).

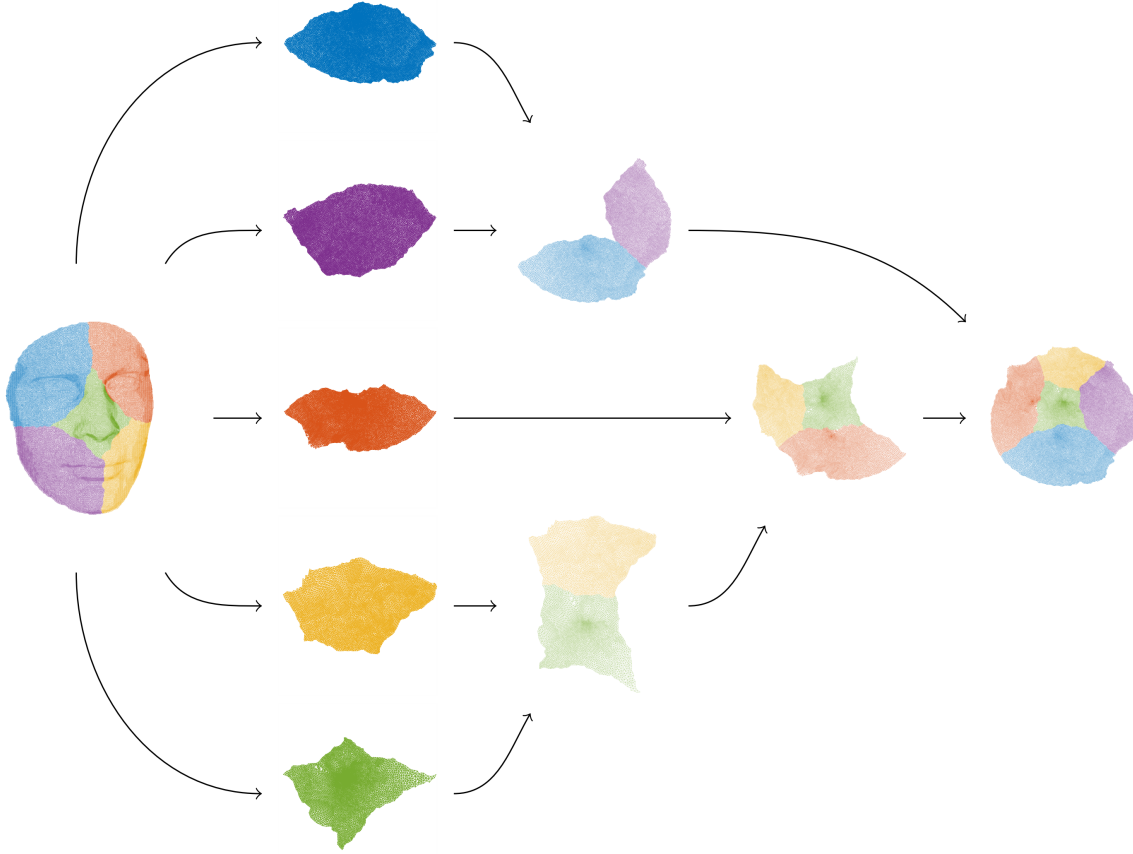


Figure 1. An overview of our proposed parallelizable global conformal parameterization (PGCP) algorithm. A simply-connected surface is first partitioned into small subdomains. The subdomains are then conformally flattened onto the plane in parallel. The flattened subdomains are subsequently stitched seamlessly by a novel partial welding technique along the common boundary arcs, thereby producing a global conformal parameterization of the surface. Note that the partial welding step only involves the boundary points of the subdomains but not their interior. The mesh structures of the subdomains are shown only for visualization purpose.

surfaces can be reduced to PDEs on the parameter domain, which are much easier to solve. Also, texture mapping on a 3D surface can be done by parameterizing it onto the 2D plane, in which textures can be easily designed. Among all surface parameterizations, one special type of parameterization is called conformal parameterization, which preserves angle and hence the local geometry of the surfaces. This is particularly important for applications such as texture mapping and remeshing, in which the angle structure plays an important role in the computation. To avoid creating computational burdens or introducing distortions, a fast and accurate method for computing conformal parameterization of surfaces is desired.

In this work, we propose a novel *parallelizable global conformal parameterization* method (abbreviated as *PGCP*) for simply-connected surfaces. Unlike the existing methods, our method uses a divide and conquer approach and exploits the nature of conformal parameterization, making the computation highly parallelizable. Figure 1 gives an overview of our proposed

method. We begin with partitioning a given surface into smaller subdomains. Then, the local conformal parameterizations of the subdomains are computed in parallel. Note that the local parameterization results are not necessarily consistent along their boundaries. Motivated by the theory of conformal welding in complex analysis, we develop a method called *partial welding* to update the boundaries of the flattened subdomains for enforcing the consistency between them. Finally, we solve the Laplace equation with the updated boundary constraints to find conformal parameterizations of the subdomains such that all of them can be glued seamlessly, ultimately forming a global conformal parameterization of the given dense surface. The bijectivity of the parameterization is guaranteed by quasi-conformal theory.

The rest of the paper is organized as follows. In Section 2, we review the related works in surface parameterization. In Section 3, we introduce the mathematical concepts involved in our work. In Section 4, we describe our proposed method for computing a global conformal parameterization of simply-connected surfaces via partial welding. Experimental results and applications are presented in Section 5 for demonstrating the effectiveness of our proposed method. We conclude our work and discuss possible future works in Section 6.

2. Related works. Surface parameterization has been widely studied in geometry processing. For an overview of the subject, readers are referred to the surveys [1–3]. It is well-known that only developable surfaces can be isometrically flattened without any distortions in area and angle. For general surfaces, it is unavoidable to introduce distortions in area or angle (or both) under parameterization. This limitation leads to two major classes of surface parameterization algorithms, namely the area-preserving parameterizations and angle-preserving (conformal) parameterizations.

Existing methods for area-preserving parameterizations include the locally authalic map [4], Lie advection [5], optimal mass transport (OMT) [6, 7], density-equalizing map (DEM) [8, 9] and stretch energy minimization (SEM) [10]. While the area elements can be preserved under area-preserving parameterizations, the angular distortion is uncontrolled. Since the angular distortion is related to the local geometry of the surfaces, it is important to minimize the angular distortion in many applications such as remeshing, texture mapping and cartography. In those cases, it is preferable to use conformal parameterization.

Existing conformal parameterization methods for simply-connected open surfaces include the discrete natural conformal parameterization (DNCP) [4]/least-square conformal mapping (LSCM) [11], Yamabe flow [12], angle-based flattening (ABF) [13–15], circle patterns [16], spectral conformal mapping (SCP) [17], conformal equivalence of triangle meshes (CETM) [18], discrete Ricci flow [19–21], quasi-conformal compositions [22–25] and conformal energy minimization (CEM) [26]. There are also some notable works on the spherical conformal parameterization of genus-0 closed surfaces, including linearization of Laplace equation [27, 28], Dirichlet energy minimization [29], folding-free global conformal mapping [30], FLASH [31] and north-south iterative scheme [32].

Note that all the above-mentioned methods compute a global conformal parameterization of a given surface by handling the entire surface directly. In case the given surface mesh is dense, the computation may be expensive. Also, in case the geometry of the input mesh is complicated, performing a global computation may lead to inaccuracy. Our work aims to overcome these problems by decomposing the input surface mesh into smaller domains

and parameterizing them in parallel. The consistency between the domains is ensured by a novel technique called partial welding, thereby forming a global conformal parameterization efficiently.

3. Mathematical background.

3.1. Harmonic map and conformal map. Following [33, 34], we introduce the following definitions of harmonic map and conformal map. Let D and Ω be simply-connected regions in \mathbb{R}^2 .

Definition 3.1 (Harmonic map). A map $\varphi : D \rightarrow \Omega$ is said to be harmonic if it minimizes the Dirichlet energy

$$(3.1) \quad E_D(\varphi) = \frac{1}{2} \int_D |\nabla \varphi|^2.$$

Definition 3.2 (Conformal map). A map $\varphi : D \rightarrow \Omega$ is said to be conformal if it satisfies

$$(3.2) \quad J \frac{\partial \varphi}{\partial x} = \frac{\partial \varphi}{\partial y},$$

where J is a rotation by $\frac{\pi}{2}$ in the tangent plane. If we write $\varphi = (\varphi_x, \varphi_y)$, the above equation can be reformulated as the following equations, known as the Cauchy-Riemann equations:

$$(3.3) \quad \begin{cases} \frac{\partial \varphi_x}{\partial x} - \frac{\partial \varphi_y}{\partial y} = 0, \\ \frac{\partial \varphi_x}{\partial y} + \frac{\partial \varphi_y}{\partial x} = 0. \end{cases}$$

To achieve conformality, we could minimize the conformal energy

$$(3.4) \quad E_C(\varphi) = \frac{1}{2} \int_D \left[\left(\frac{\partial \varphi_x}{\partial x} - \frac{\partial \varphi_y}{\partial y} \right)^2 + \left(\frac{\partial \varphi_x}{\partial y} + \frac{\partial \varphi_y}{\partial x} \right)^2 \right].$$

As shown by Hutchinson [33], if we define the area $A(\varphi)$ by

$$(3.5) \quad A(\varphi) = \int_D \left\| \frac{\partial \varphi}{\partial x} \times \frac{\partial \varphi}{\partial y} \right\|,$$

then the conformal energy can be expressed in terms of the Dirichlet energy and area:

$$(3.6) \quad E_C(\varphi) = E_D(\varphi) - A(\varphi).$$

Since the conformal energy is nonnegative, it follows that the Dirichlet energy is always bounded below by the area. In particular, the equality holds if and only if φ is conformal.

Moreover, given the area term $A(\varphi)$, minimizing the conformal energy is equivalent to minimizing the Dirichlet energy. Note that the area depends on how φ maps the boundary. In other words, given a “good” boundary condition, a conformal map can be obtained by simply finding the harmonic map under the given boundary condition.

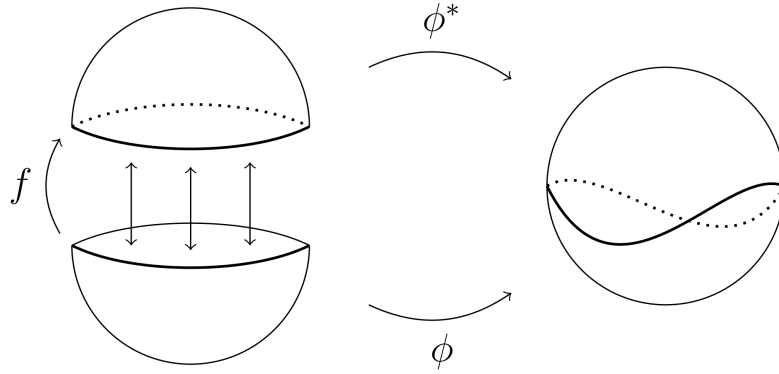


Figure 2. An illustration of the closed welding problem. The entire boundaries of the two parts are glued consistently.

3.2. Möbius transformation. A special type of conformal maps on the extended complex plane $\overline{\mathbb{C}}$ are the *Möbius transformations*, also known as the *linear fractional transformations*:

Definition 3.3 (Möbius transformation). A function $f : \overline{\mathbb{C}} \rightarrow \overline{\mathbb{C}}$ is said to be Möbius transformation if it is of the form

$$(3.7) \quad f(z) = \frac{az + b}{cz + d},$$

where a, b, c, d are complex numbers with $ab - bc \neq 0$.

Given two sets of distinct points $\{z_1, z_2, z_3\}$ and $\{w_1, w_2, w_3\}$, there exists a unique Möbius transformation satisfying $f(z_i) = w_i$, $i = 1, 2, 3$. Therefore, Möbius transformations provides us with a simple way of fixing three points conformally.

3.3. Conformal welding. *Conformal welding*, also known as *sewing* or simply *welding*, is a problem in complex analysis which concerns with gluing two surfaces in a conformal way so that they fit together consistently according to certain correspondence.

Given a diffeomorphism f from a curve (e.g. the unit circle) to itself, we want to find two Jordan domains $D, \Omega \subset \overline{\mathbb{C}}$ and two conformal maps $\phi : D \rightarrow \Omega$ and $\phi^* : D^* \rightarrow \Omega^*$ such that $\phi = \phi^* \circ f$ on the curve [35]. Here, D^* and Ω^* are the exterior of D and Ω respectively. Since $\overline{\mathbb{C}} \cong \mathbb{S}^2$, the two domains D, Ω can be regarded as two disk-like surfaces on \mathbb{S}^2 . Intuitively, given a correspondence between the boundaries of the two surfaces, the problem of conformal welding is to find two conformal deformations such that the surfaces are stitched together seamlessly (see Figure 2). We refer this classical welding problem as a *closed* welding problem.

For a general homeomorphism f , the closed welding problem may not have any solution. However, if f satisfies certain conditions, the problem is solvable. We introduce the concept of *quasisymmetric function* below:

Definition 3.4 (Quasisymmetric function [36]). Let f be a continuous, strictly increasing function defined on an interval I of the x -axis. We call f k -quasisymmetric (or simply

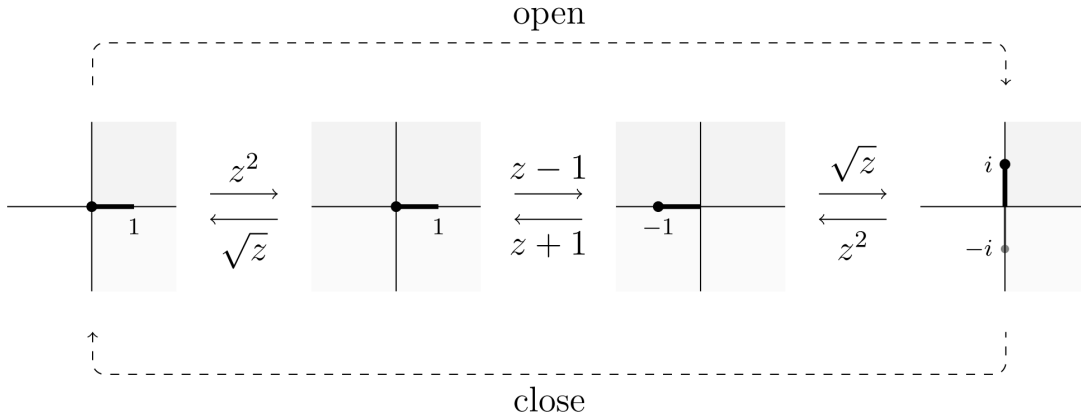


Figure 3. An illustration of the opening map and the closing map. The top part shows the processes involved in the opening map $z \mapsto \sqrt{z^2 - 1}$, which ultimately map $\{0, 1\}$ to $\{i, 0\}$ or $\{-i, 0\}$, depending on the choice of branching. The closing map $z \mapsto \sqrt{z^2 + 1}$ reverses the processes such that $\pm i$ will be mapped back to 0.

quasisymmetric) on I if there exists a positive constant k such that

$$(3.8) \quad \frac{1}{k} \leq \frac{f(x+t) - f(x)}{f(x) - f(x-t)} \leq k$$

for all $x, x-t, x+t \in I$ with $t > 0$.

One can show that the closed welding problem is solvable if f is a quasisymmetric function from the real axis to itself:

Theorem 3.5 (Sewing theorem [36]). Let f be a quasisymmetric function on the real axis. Then the upper and lower half-planes can be mapped conformally onto disjoint Jordan domains D, Ω by two maps ϕ, ϕ^* , with $\phi(x) = \phi^*(f(x))$ for all $x \in \mathbb{R}$.

The theorem was first proven by Pfluger based on the existence of solutions to the Beltrami equation [37]. Another way to prove the result is to use some approximation techniques on the quasisymmetric function [36].

3.4. Geodesic algorithm. A conformal mapping method called the *zipper algorithm* was proposed independently by Kühnau [38] and Marshall and Morrow [39] in the 1980s. In particular, Marshall and Rohde [40] proved the convergence of a variant of it called the *geodesic algorithm*. The geodesic algorithm computes a conformal map from a region in the complex plane to the upper half-plane \mathbb{H} . Below, we briefly describe the geodesic algorithm.

The key ingredients of the geodesic algorithm are two maps: the *opening map* $z \mapsto \sqrt{z^2 - 1}$ and the *closing map* $z \mapsto \sqrt{z^2 + 1}$. Intuitively, they are operations analogous to opening and closing a slit, behaving like a zipper (see Figure 3). Suppose we have a simple closed region Ω , and a sequence of boundary points $\{z_0, z_1, \dots, z_k\}$ on $\partial\Omega$. To initiate the process, define a map $g_1 : \Omega \rightarrow \mathbb{C}$ by

$$(3.9) \quad g_1(z) = \sqrt{\frac{z - z_1}{z - z_0}},$$

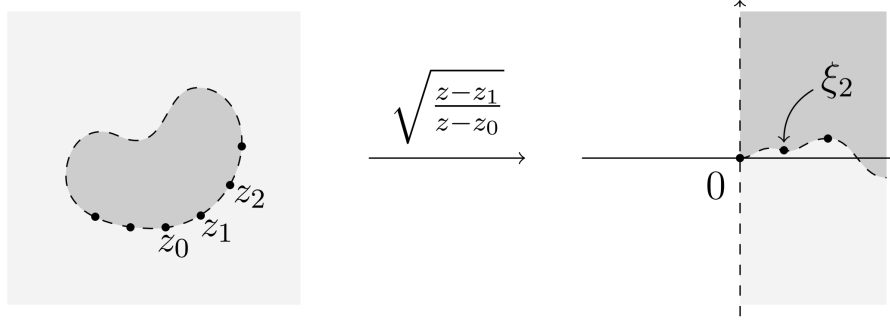


Figure 4. In the geodesic algorithm, the first map g_1 maps z_0 to ∞ and z_1 to 0 . It initiates the process so that we can apply the opening maps for the remaining data points.

164 assuming the branching $(-1)^{1/2} = i$. This maps Ω to the right half-plane. In particular, the
 165 line segment between z_0 and z_1 is mapped onto the imaginary axis, with z_0 mapped to ∞ and
 166 z_1 mapped to 0 (see Figure 4).

167 Analogously, one can define a map g_2 such that the line segment between $g_1(z_1) = 0$ and
 168 $\xi_2 := g_1(z_2)$ is mapped to the imaginary axis, while the remaining points are still in the right
 169 half-plane. By further repeating the above process, all the boundary points can be pushed
 170 onto the imaginary axis one by one. More explicitly, suppose the point z_j has already been
 171 transformed to the position ξ_j after applying the opening maps g_1, g_2, \dots, g_{j-1} , i.e.

$$172 \quad (3.10) \quad \xi_j = g_{j-1} \circ g_{j-2} \circ \dots \circ g_1(z_j).$$

173 Consider a Möbius transformation

$$174 \quad (3.11) \quad L_{\xi_j}(z) := \frac{\frac{\operatorname{Re}(\xi_j)}{|\xi_j|^2} z}{1 + \frac{\operatorname{Im}(\xi_j)}{|\xi_j|^2} zi}.$$

175 It can be easily checked that L_{ξ_j} maps $\{0, \xi_j, \rho\}$ to $\{0, 1, \infty\}$, where ρ is a point on the imaginary
 176 axis at which the orthogonal circular arc from 0 to ξ_k extends to. Now, as $L_{\xi_j}(\xi_j) = 1$, we
 177 can map the segment between $L_{\xi_j}(g_{j-1} \circ g_{j-2} \circ \dots \circ g_1(z_{j-1})) = 0$ and $L_{\xi_j}(\xi_j) = 1$ onto the
 178 imaginary axis as illustrated in Figure 3. Define g_j as the composition of L_{ξ_j} with the opening
 179 map $f(z) = \sqrt{z^2 - 1}$:

$$180 \quad (3.12) \quad g_j(z) := \sqrt{L_{\xi_j}(z)^2 - 1}.$$

181 Note that $g_j((g_{j-1} \circ \dots \circ g_1(z_{j-1}))) = \sqrt{0 - 1} = i$ (assuming the branching $(-1)^{1/2} = i$),
 182 $g_j(\xi_j) = \sqrt{L_{\xi_j}(\xi_j)^2 - 1} = \sqrt{1 - 1} = 0$, and the entire region will remain in the right half-plane.
 183 Therefore, the requirements for g_j are satisfied.

184 After obtaining the maps g_1, g_2, \dots, g_k such that $g_k \circ g_{k-1} \circ \dots \circ g_1$ maps all boundary
 185 points $\{z_0, z_1, \dots, z_k\}$ onto the imaginary axis, define a final map

$$186 \quad (3.13) \quad g_{k+1}(z) = \left(\frac{z}{1 - \frac{z}{g_k \circ g_{k-1} \circ \dots \circ g_1(z_0)}} \right)^2.$$

g_{k+1} maps the transformed region $g_k \circ g_{k-1} \circ \cdots \circ g_1(\Omega)$ onto the upper half-plane \mathbb{H} . Since all the above maps are analytic and, in particular, a square and a square root map are applied in each step, the composition map $(g_{k+1} \circ g_k \circ \cdots \circ g_1) : \Omega \rightarrow \mathbb{H}$ is conformal.

3.5. Quasi-conformal map. Quasi-conformal map is an extension of conformal map in the sense that it allows for bounded conformal distortion. Intuitively, conformal maps map infinitesimal circles to infinitesimal circles, while quasi-conformal maps map infinitesimal circles to infinitesimal ellipses with bounded eccentricity. The formal definition of quasi-conformal map is given below.

Definition 3.6 (Quasi-conformal map [41]). A map $\varphi : D \rightarrow \Omega$ is said to be quasi-conformal if it satisfies the Beltrami equation

$$(3.14) \quad \frac{\partial \varphi}{\partial \bar{z}} = \mu_\varphi(z) \frac{\partial \varphi}{\partial z}$$

for some complex-valued function μ_φ with $\|\mu_\varphi\|_\infty < 1$. μ_φ is said to be the Beltrami coefficient of φ .

The Beltrami coefficient μ_φ captures the conformal distortion of φ . In particular, if $\mu_\varphi = 0$, then the Beltrami equation becomes the Cauchy-Riemann equations and hence φ is conformal. Also, the Jacobian J_φ of φ is given by

$$(3.15) \quad J_\varphi = \left| \frac{\partial \varphi}{\partial z} \right|^2 (1 - |\mu_\varphi|^2).$$

Therefore, a map is folding-free if and only if its Beltrami coefficient is with sup norm less than 1.

Moreover, one can correct the conformal distortion and non-bijectivity of a map by composing it with another map. If $\varphi_1 : \mathbb{C} \rightarrow \mathbb{C}$ and $\varphi_2 : \mathbb{C} \rightarrow \mathbb{C}$ are two maps with Beltrami coefficients μ_{φ_1} and μ_{φ_2} , then $\varphi_2 \circ \varphi_1$ is a quasi-conformal map with Beltrami coefficient

$$(3.16) \quad \mu_{\varphi_2 \circ \varphi_1}(z) = \frac{\mu_{\varphi_1}(z) + \frac{\overline{\varphi_{1z}}}{\varphi_{1z}} \mu_{\varphi_2}(\varphi_1(z))}{1 + \frac{\overline{\varphi_{1z}}}{\varphi_{1z}} \mu_{\varphi_1}(z) \mu_{\varphi_2}(\varphi_1(z))}.$$

In particular, if $\mu_{\varphi_2} = \mu_{\varphi_1}^{-1}$, then $\mu_{\varphi_2 \circ \varphi_1} = 0$ and hence the composition map $\varphi_2 \circ \varphi_1$ is conformal and folding-free. This idea of quasi-conformal composition has been used in [24, 25, 31], and the details of the theory and computation of it can be found therein.

4. Proposed method. Let \mathcal{S} be a simply-connected surface in \mathbb{R}^3 , with a triangle mesh representation $(\mathcal{V}, \mathcal{F})$ where \mathcal{V} is the vertex set and \mathcal{F} is the face set. Our goal is to compute a global conformal parameterization of $\mathcal{S} = (\mathcal{V}, \mathcal{F})$ in an efficient and accurate way.

4.1. Surface partition. The first step is to partition \mathcal{S} into submeshes based on a prescribed set of edges $\tilde{\mathcal{E}}$. More specifically, denote the edge set of \mathcal{S} by \mathcal{E} , and the set of boundary edges of \mathcal{S} by \mathcal{E}_{bdy} . Consider the set $E = \mathcal{E} \setminus (\tilde{\mathcal{E}} \cup \mathcal{E}_{\text{bdy}})$. We construct a graph G using E and find all connected components in G . Suppose there are K connected components in G , where each of them consists of a sub-face set \mathcal{F}_i , $i = 1, \dots, K$. By tracking all vertices that are contained

in \mathcal{F}_i , we obtain a sub-vertex set \mathcal{V}_i . In other words, we have obtained K simply-connected open submeshes $\mathcal{S}_1 = (\mathcal{V}_1, \mathcal{F}_1), \mathcal{S}_2 = (\mathcal{V}_2, \mathcal{F}_2), \dots, \mathcal{S}_K = (\mathcal{V}_K, \mathcal{F}_K)$ that satisfy the following properties:

(i) The union of the vertex sets of all submeshes is exactly \mathcal{V} :

$$(4.1) \quad \bigcup_{i=1}^K \mathcal{V}_i = \mathcal{V}.$$

(ii) The union of the face sets of all submeshes is exactly \mathcal{F} :

$$(4.2) \quad \bigcup_{i=1}^K \mathcal{F}_i = \mathcal{F}.$$

(iii) The intersection of any two different sub-vertex sets is the intersection of the boundary sets, which is either an empty set or a boundary segment:

$$(4.3) \quad \mathcal{V}_i \cap \mathcal{V}_j = \partial \mathcal{S}_i \cap \partial \mathcal{S}_j \text{ for all } i, j.$$

(iv) The intersection of any two different sub-face sets is empty:

$$(4.4) \quad \mathcal{F}_i \cap \mathcal{F}_j = \emptyset \text{ for all } i \neq j.$$

4.2. Local conformal parameterization of submeshes. The next step is to compute a conformal parameterization of every \mathcal{S}_i . To find a conformal parameterization $\varphi_i : \mathcal{S}_i \rightarrow \mathbb{R}^2$, the DNCP method [4] is used. In short, DNCP minimizes the Dirichlet energy $E_D(\varphi)$ and maximizes the area $A(\varphi)$, based on the fact that the Dirichlet energy is bounded below by the area and conformality is attained when equality holds. We briefly describe the method below.

Let $\mathcal{V}_i = \{v_{i_1}, v_{i_2}, \dots, v_{i_{n_i}}\}$ be the vertices in \mathcal{S}_i , and $\varphi_i : \mathcal{S}_i \rightarrow \mathbb{R}^2$ be a flattening map. Denote $\mathbf{u} = [\mathbf{u}_1, \mathbf{u}_2, \dots, \mathbf{u}_{n_i}]^t = [\varphi_i(v_{i_1}), \varphi_i(v_{i_2}), \dots, \varphi_i(v_{i_{n_i}})]^t$. The Dirichlet energy is discretized using the cotangent formula [34]:

$$(4.5) \quad E_D(\mathbf{u}) = \frac{1}{2} \sum_{(v_{i_p}, v_{i_q}) \text{ adjacent}} (\cot \alpha_{pq} + \cot \beta_{pq}) |\mathbf{u}_p - \mathbf{u}_q|^2 = \mathbf{u}^t L^{\cotan} \mathbf{u},$$

where α_{pq}, β_{pq} are the two angles opposite to the edge $[v_{i_p}, v_{i_q}]$ in \mathcal{S}_i , and L^{\cotan} is a $|\mathcal{V}_i| \times |\mathcal{V}_i|$ sparse symmetric positive definite matrix also known as the cotangent Laplacian:

$$(4.6) \quad L_{p,q}^{\cotan} = \begin{cases} \frac{1}{2}(\cot \alpha_{pq} + \cot \beta_{pq}) & \text{if } (v_{i_p}, v_{i_q}) \text{ are adjacent,} \\ -\sum_{r \neq p} L_{p,r}^{\cotan} & \text{if } p = q, \\ 0 & \text{otherwise.} \end{cases}$$

The area is discretized using the boundary vertices of \mathcal{S}_i :

$$(4.7) \quad A(\mathbf{u}) = \frac{1}{2} \sum_{[v_{i_p}, v_{i_q}] \subset \partial \mathcal{S}_i} (x_p y_q - y_p x_q) = (\mathbf{x}^t \quad \mathbf{y}^t) M^{\text{area}} \begin{pmatrix} \mathbf{x} \\ \mathbf{y} \end{pmatrix},$$

where $\mathbf{u}_j = (x_j, y_j)$ for all j , $\mathbf{x} = [x_1, x_2, \dots, x_{n_i}]^t$ is the collection of all x -coordinates of \mathbf{u} , $\mathbf{y} = [y_1, y_2, \dots, y_{n_i}]^t$ is the collection of all y -coordinates of \mathbf{u} , and M^{area} is a $2|\mathcal{V}_i| \times 2|\mathcal{V}_i|$ sparse symmetric matrix. More explicitly, if $[v_{i_p}, v_{i_q}] \subset \partial\mathcal{S}_i$, we have

$$(4.8) \quad M_{p,q+|\mathcal{V}_i|}^{\text{area}} = M_{q+|\mathcal{V}_i|,p}^{\text{area}} = 1 \quad \text{and} \quad M_{q,p+|\mathcal{V}_i|}^{\text{area}} = M_{p+|\mathcal{V}_i|,q}^{\text{area}} = -1.$$

DNCP minimizes the discrete conformal energy

$$(4.9) \quad E_C(\mathbf{u}) = E_D(\mathbf{u}) - A(\mathbf{u})$$

subject to the prescribed positions of two boundary vertices that remove the freedom of rigid motion and scaling. It suffices to solve a $2|\mathcal{V}_i| \times 2|\mathcal{V}_i|$ sparse linear system

$$(4.10) \quad \left(\begin{pmatrix} L^{\cotan} & 0 \\ 0 & L^{\cotan} \end{pmatrix} - M^{\text{area}} \right) \begin{pmatrix} \mathbf{x} \\ \mathbf{y} \end{pmatrix} = 0$$

subject to four boundary constraints (two in \mathbf{x} and two in \mathbf{y} for the two pinned boundary vertices). The resultant map φ_i satisfying $\varphi_i(\mathcal{V}_i) = \mathbf{u} = [\mathbf{x}, \mathbf{y}]$ is the desired conformal parameterization of \mathcal{S}_i .

DNCP is suitable for our framework since it is a free-boundary linear method. As discussed above, obtaining each φ_i only requires solving a $2|\mathcal{V}_i| \times 2|\mathcal{V}_i|$ sparse matrix equation, which is highly efficient. Also, the free-boundary condition ensures that no additional conformal distortion will be introduced at the boundaries. This is particularly important in our subsequent welding step.

It is noteworthy that the parameterization of each submesh is independent, and hence this step of computing local conformal parameterizations is highly parallelizable.

4.3. Partial welding. Note that the local parameterizations we obtained via DNCP are not necessarily consistent along the boundaries. Therefore, we need a step for gluing the boundaries of them consistently. To preserve the conformality of the parameterization, the gluing step should be conformal. This problem of gluing subdomains is different from the closed welding problem introduced in Section 3. More explicitly, the closed welding problem considers gluing the entire boundaries of two domains, while in general only a portion of the boundaries of two neighboring subdomains in our case should be glued. In other words, the problem that we need to tackle is a *partial* welding problem that involves gluing two subdomains along only a pair of boundary arcs.

Below, we first rigorously derive a theoretical construction for solving the partial welding problem. Then, we devise an efficient algorithm for solving it.

4.3.1. Theoretical construction. We formulate the problem mathematically. Given two Jordan regions $A, B \subset \overline{\mathbb{C}}$, let $\gamma_A \subset \partial A$ and $\gamma_B \subset \partial B$ be some arcs of the boundaries of A and B respectively. Suppose we have a correspondence function $f : \gamma_A \rightarrow \gamma_B$ that relates points on γ_A and points on γ_B . The partial welding problem is to find two conformal maps $\Phi_A : A \rightarrow A'$ and $\Phi_B : B \rightarrow B'$, with A' and B' being disjoint, such that

$$(4.11) \quad \Phi_A(\gamma_A) = (\Phi_B \circ f)(\gamma_A).$$

Recall that the closed welding problem is solvable for quasisymmetric function on the real axis. For the partial welding problem, we make use of the following lemma.

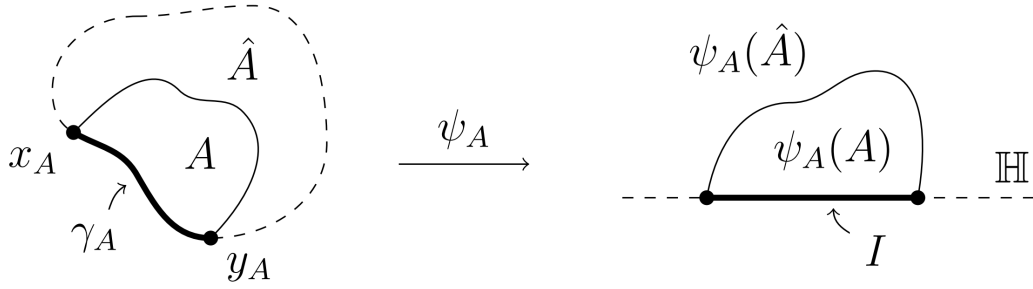


Figure 5. Construction of \hat{A} and the mapping from it to the upper half plane.

285 **Lemma 4.1** (Lehto and Virtanen [36]). *Every function f which is k -quasisymmetric on an*
 286 *interval $I = [a, b]$ can be extended to a k -quasisymmetric function on the entire x -axis, where*
 287 *the constant k is less than a number depending only on k .*

288 To make use of the above lemma, we further suppose that \hat{A} is a larger domain that
 289 contains A while sharing the boundary segment γ_A , i.e. $A \subset \hat{A}$ and $\gamma_A \subset \partial\hat{A}$ (see Figure 5
 290 left). Denote the endpoints of γ_A by x_A and y_A . Similarly, let \hat{B} be a domain such that $B \subset \hat{B}$
 291 and $\gamma_B \subset \partial\hat{B}$, and denote the endpoints of γ_B by x_B and y_B .

292 By the Riemann mapping theorem, \hat{A} and \hat{B} can be mapped to the upper and lower
 293 half plane respectively by some conformal maps ψ_A and ψ_B . Now, we fix x_A and y_A at the
 294 endpoints of some interval I on the x -axis. For simplicity, we take $I = [-1, 1]$ and fix x_A and
 295 y_A at -1 and 1 respectively (see Figure 5 right). Similarly, we fix x_B at -1 and y_B at 1 . The
 296 homeomorphic extensions to the closures define a map $g : I \rightarrow I$ by $g = \psi_B \circ f \circ \psi_A^{-1}$. In
 297 other words, we have $f = \psi_B^{-1} \circ g \circ \psi_A$ by construction. Assuming that g is a quasisymmetric
 298 function from I to itself, we get a quasisymmetric extension $\hat{g} : \mathbb{R} \rightarrow \mathbb{R}$ of g using Lemma 4.1.

299 Then, we apply Theorem 3.5 with this \hat{g} , which gives us two conformal maps $\phi_A : \hat{A} \rightarrow \hat{A}'$
 300 and $\phi_B : \hat{B} \rightarrow \hat{B}'$ with \hat{A}' and \hat{B}' being disjoint, such that the boundary values satisfy
 301 $\phi_A(x) = \phi_B(\hat{g}(x))$ for all $x \in \mathbb{R}$. In particular, $\phi_A(x) = \phi_B(g(x))$ for all $x \in I$. Figure 6 shows
 302 an illustration of the construction.

303 Since the composition of conformal maps is conformal, we have constructed two conformal
 304 maps $\Phi_A = \phi_A \circ \psi_A$ and $\Phi_B = \phi_B \circ \psi_B$, which respectively map A to some $A' \subset \hat{A}'$ and B to
 305 some $B' \subset \hat{B}'$. Note that $f = \psi_B^{-1} \circ g \circ \psi_A$ when we restrict f on γ_A . Also,

$$(4.12) \quad (\phi_B \circ \psi_B) \circ f = \phi_B \circ \psi_B \circ \psi_B^{-1} \circ g \circ \psi_A = \phi_B \circ g \circ \psi_A = \phi_A \circ \psi_A,$$

307 where the last equality follows from Theorem 3.5. This solves the partial welding problem.

308 **4.3.2. Algorithmic construction.** The theoretical construction above provides us with
 309 a continuous approach for solving the partial welding problem. We proceed to develop
 310 an algorithm to solve the problem over discrete boundary data points. Suppose we have
 311 two sequences of boundary points $\partial A = \{a_0, \dots, a_k, \dots, a_m\}$ and $\partial B = \{b_0, \dots, b_k, \dots, b_n\}$,
 312 where a_j corresponds to b_j (i.e. a_j should be glued with b_j) for $j = 0, \dots, k$. This gives
 313 a correspondence function $f : \gamma_A \subset \partial A \rightarrow \gamma_B \subset \partial B$, where $\gamma_A = \{a_0, \dots, a_k\}$ and $\gamma_B =$

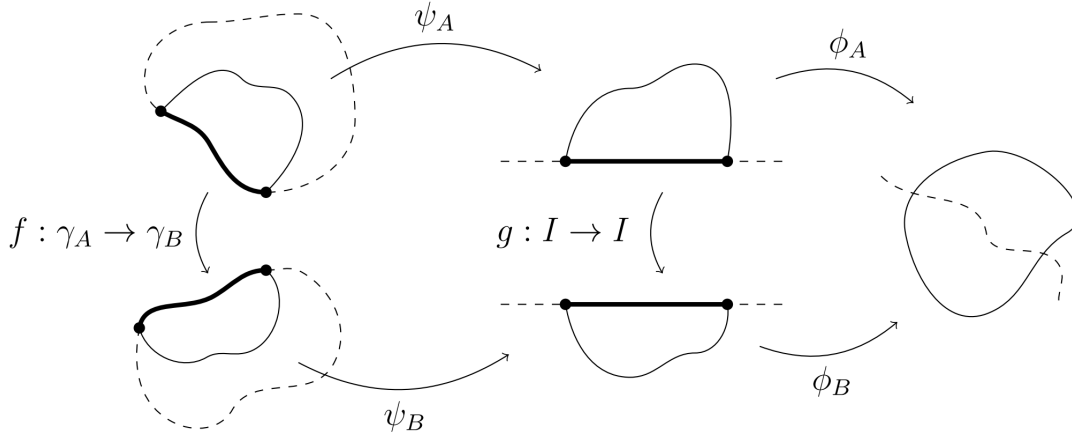


Figure 6. The construction of conformal maps for solving the partial welding problem.

$\{b_0, \dots, b_k\}$, with $f(a_j) = b_j$ for $j = 0, \dots, k$. Our goal is to construct the two maps Φ_A, Φ_B for gluing the two boundary curves conformally along the corresponding points. As discussed in the theoretical construction, a good way for the construction of Φ_A, Φ_B is to make use of two maps ψ_A, ψ_B that map A, B onto the upper and lower half-planes respectively. We propose an algorithm that makes use of a variant of the geodesic algorithm [40].

We begin with designing an algorithm that maps a sequence of boundary points to a standard shape. The algorithm is based on a key observation that the geodesic algorithm can be paused halfway. Suppose we have a sequence of boundary points $\{z_0, \dots, z_k, \dots, z_n\}$. Consider applying the first k maps g_1, g_2, \dots, g_k in the geodesic algorithm on $\{z_0, \dots, z_k, \dots, z_n\}$ with branching $(-1)^{1/2} = i$. The composition $g_k \circ \dots \circ g_1$ maps the first $k+1$ points $\{z_0, \dots, z_k\}$ onto the imaginary axis, with z_k mapped to 0, while the remaining boundary data points $\{z_{k+1}, \dots, z_n\}$ are all mapped onto the right half-plane. Note that each of g_2, \dots, g_k is a composition of a Möbius transformation, a square map and a square root map. Therefore, they are all conformal.

Now, instead of the final map (3.13) in the geodesic algorithm, we apply the following Möbius transformation:

$$(4.13) \quad g_{k+1}(z) = \frac{z}{1 - \frac{z}{g_k \circ g_{k-1} \circ \dots \circ g_1(z_0)}}.$$

It is easy to check that $g_{k+1}(g_k \circ \dots \circ g_1(z_0)) = \infty$ and $g(0) = 0$. In other words, the new composition $g_{k+1} \circ g_k \circ \dots \circ g_1$ maps the first data point z_0 to ∞ and the $(k+1)$ -th data point z_k to 0. Note that the first $k+1$ data points are on the upper half of the imaginary axis, and the remaining boundary data points $\{z_{k+1}, \dots, z_n\}$ are on the right half-plane. We call such a half-opened (i.e. half-unzipped) shape an *intermediate form*. Note that by using another branching $(-1)^{1/2} = -i$ throughout the maps above, we have an alternative way to transform a sequence of boundary data points onto the right half-plane, with the first $k+1$ data points mapped onto the lower half of the imaginary axis. Algorithm 1 summarizes the proposed intermediate form transformation procedure.

Algorithm 1: Intermediate form transformation

Input: A sequence of boundary points $\{z_0, \dots, z_k, \dots, z_n\}$ and a choice of branching.

Output: A sequence of transformed boundary points $\{Z_0, \dots, Z_k, \dots, Z_n\}$, where Z_0, \dots, Z_k are on the imaginary axis.

- 1 Set $g_1(z) = \sqrt{\frac{z-z_1}{z-z_0}}$ with the choice of branching;
- 2 **for** $j = 2, \dots, k$ **do**
- 3 Compute $\xi_j = (g_{j-1} \circ \dots \circ g_1)(z_j)$;
- 4 Set $g_j(z) = \sqrt{L_{\xi_j}(z)^2 - 1}$ with the choice of branching, where $L_{\xi_j}(z) := \frac{\frac{\operatorname{Re}(\xi_j)}{|\xi_j|^2}z}{1 + \frac{\operatorname{Im}(\xi_j)}{|\xi_j|^2}zi}$;
- 5 Set $g_{k+1}(z) = \frac{z}{1 - \frac{z}{g_k \circ g_{k-1} \circ \dots \circ g_1(z_0)}}$;
- 6 Compute $P_l = (g_{k+1} \circ \dots \circ g_1)(z_l)$ for $l = 0, \dots, k, \dots, n$;

Coming back to the problem of aligning the two sequences of boundary points $\partial A = \{a_0, \dots, a_k, \dots, a_m\}$ and $\partial B = \{b_0, \dots, b_k, \dots, b_n\}$, we define auxiliary data points $a_{m+1} = b_{n+1} = 0, a_{m+2} = b_{n+2} = \infty$ to keep track of the transformation. Now, using Algorithm 1 with two different choices of branching $(-1)^{1/2} = i$ and $(-1)^{1/2} = -i$, we map $\{a_0, \dots, a_k, \dots, a_m, a_{m+1}, a_{m+2}\}$ and $\{b_0, \dots, b_k, \dots, b_n, b_{n+1}, b_{n+2}\}$ onto the right half-plane. Denote the transformed data points by $\{A_0, \dots, A_k, \dots, A_m, A_{m+1}, A_{m+2}\}$ and $\{B_0, \dots, B_k, \dots, B_n, B_{n+1}, B_{n+2}\}$. Note that $\{A_0, \dots, A_k\}$ are all on the upper half of the imaginary axis with $A_0 = \infty$ and $A_k = 0$, while $\{B_0, \dots, B_k\}$ are all on the lower half of the imaginary axis with $B_0 = \infty$ and $B_k = 0$. The next step is to align A_j with B_j for all $j = 0, \dots, k$ conformally, such that the two boundary curves ∂A and ∂B are welded based on the partial correspondence between γ_A and γ_B .

Suppose $\alpha = ai$ and $\beta = bi$ are two corresponding points originally on γ_A and γ_B under the intermediate form transformations, where $a > 0 > b$. A Möbius transformation that takes $\{\alpha, 0, \beta\}$ to $\{i, 0, -i\}$ is explicitly given by

$$(4.14) \quad T_\beta^\alpha(z) = \frac{z}{\frac{-2ab}{a-b} - \frac{a+b}{a-b}zi}.$$

This transformation provides us with a simple way to align each pair of corresponding points. Note that $A_k = 0 = B_k$ is automatically aligned, and so we start with aligning A_{k-1} and B_{k-1} . Applying the Möbius transformation $T_{B_{k-1}}^{A_{k-1}}$ onto the two sets of boundary data points, we map A_{k-1} to i and B_{k-1} to $-i$. Then, we compose the map with the closing map $z \mapsto \sqrt{z^2 + 1}$ so that i and $-i$ are both mapped to 0. More explicitly, we define

$$(4.15) \quad h_{k-1}(z) := \sqrt{T_{B_{k-1}}^{A_{k-1}}(z)^2 + 1}$$

and apply it to all data points. The branching for the computation of each point is determined using the previous choice in the intermediate form transformation. Then, we repeat the above

process for $j = k - 2, \dots, 1$ by defining

$$(4.16) \quad h_j(z) := \sqrt{T_{\beta_j}^{\alpha_j}(z)^2 + 1},$$

where

$$(4.17) \quad \alpha_j = (h_{j+1} \circ \dots \circ h_{k-1})(A_j)$$

and

$$(4.18) \quad \beta_j = (h_{j+1} \circ \dots \circ h_{k-1})(B_j).$$

Now, all pairs of corresponding points $(A_1, B_1), \dots, (A_k, B_k)$ have been consistently aligned under the composition map $h_1 \circ h_2 \circ \dots \circ h_{k-1}$. The first pair of corresponding points $A_0 = \infty = B_0$ are also automatically aligned. Note that each of h_{k-1}, \dots, h_1 is a composition of a Möbius transformation, a square map and a square root map. Hence, they are all conformal.

Then, we define a closing map h_0 similar to (3.13) in the geodesic algorithm:

$$(4.19) \quad h_0(z) := \left(\frac{z}{1 - \frac{z}{(h_1 \circ \dots \circ h_k)(\infty)}} \right)^2.$$

Note that h_0 maps all points onto the upper half plane \mathbb{H} , with 0 mapped to 0 and $(h_1 \circ \dots \circ h_{k-1})(\infty)$ (i.e. $(h_1 \circ \dots \circ h_{k-1})(A_0)$) mapped to ∞ . We obtain the transformed data points

$$(4.20) \quad \tilde{a}_l = (h_0 \circ \dots \circ h_{k-1})(A_l)$$

for $l = 0, \dots, m + 2$ with branching $(-1)^{1/2} = i$, and

$$(4.21) \quad \tilde{b}_l = (h_0 \circ \dots \circ h_{k-1})(B_l)$$

for $l = 0, \dots, n + 2$ with branching $(-1)^{1/2} = -i$.

Considering the entire composition $h_0 \circ h_1 \circ \dots \circ h_{k-1} \circ g_{k+1} \circ \dots \circ g_1$ starting from the beginning to here, it can be observed that $g_2, \dots, g_{k+1}, h_{k-1}, \dots, h_2$ are all conformal, while g_2 is a square root map and h_0 is a square map. Therefore, the entire composition is conformal. In other words, we have conformally transformed the two sequences of boundary data points $\{a_0, \dots, a_k, \dots, a_m\}$ and $\{b_0, \dots, b_k, \dots, b_n\}$ into $\{\tilde{a}_0, \dots, \tilde{a}_k, \dots, \tilde{a}_m\}$ and $\{\tilde{b}_0, \dots, \tilde{b}_k, \dots, \tilde{b}_n\}$ such that the partial correspondence between them is satisfied, i.e. $\tilde{a}_j = \tilde{b}_j$ for $j = 0, \dots, k$.

Finally, we perform a normalization by tracking the transformation of the auxiliary data points $a_{m+1} = b_{n+1} = 0, a_{m+2} = b_{n+2} = \infty$. More explicitly, we apply a Möbius transformation T that takes $\{\tilde{a}_{m+1}, \tilde{b}_{n+1}, \frac{1}{2}(\tilde{a}_{m+2} + \tilde{a}_{n+2})\}$ to $\{-1, 1, \infty\}$ on all the transformed points. This regularizes the transformation and prevents the boundary data points from being mapped far away. Note that Möbius transformations are conformal and hence the conformality of the composition map is preserved. This completes the process of gluing two boundary curves based on a partial correspondence between them. Algorithm 2 summarizes the proposed partial welding algorithm.

Algorithm 2: Partial welding

Input: Two sequences of boundary data points $\{a_0, \dots, a_k, \dots, a_m\}$ and $\{b_0, \dots, b_k, \dots, b_n\}$, where a_j should be glued with b_j for $j = 0, \dots, k$.

Output: Conformally transformed data points $\{\tilde{a}_0, \dots, \tilde{a}_k, \dots, \tilde{a}_m\}$ and $\{\tilde{b}_0, \dots, \tilde{b}_k, \dots, \tilde{b}_n\}$ such that $\tilde{a}_j = \tilde{b}_j$ for $j = 0, \dots, k$.

- 1 Define auxiliary points $a_{m+1} = b_{n+1} = 0, a_{m+2} = b_{n+2} = \infty$;
- 2 Apply Algorithm 1 on $\{a_0, \dots, a_k, \dots, a_m, a_{m+1}, a_{m+2}\}$ with branching $(-1)^{1/2} = i$ and obtain the transformed boundary data points $\{A_0, \dots, A_k, \dots, A_m, A_{m+1}, A_{m+2}\}$;
- 3 Apply Algorithm 1 on $\{b_0, \dots, b_k, \dots, b_n, b_{n+1}, b_{n+2}\}$ with branching $(-1)^{1/2} = -i$ and obtain the transformed boundary data points $\{B_0, \dots, B_k, \dots, B_n, B_{n+1}, B_{n+2}\}$;
- 4 Set $h_{k-1}(z) := \sqrt{T_{B_{k-1}}^{A_{k-1}}(z)^2 + 1}$;
- 5 **for** $j = k-2, \dots, 1$ **do**
- 6 Compute $\alpha_j = (h_{j+1} \circ \dots \circ h_{k-1})(A_j)$ with branching $(-1)^{1/2} = i$;
- 7 Compute $\beta_j = (h_{j+1} \circ \dots \circ h_{k-1})(B_j)$ with branching $(-1)^{1/2} = -i$;
- 8 Set $h_j(z) := \sqrt{T_{\beta_j}^{\alpha_j}(z)^2 + 1}$;
- 9 Set $h_0(z) := \left(\frac{z}{1 - \frac{z}{(h_1 \circ \dots \circ h_{k-1})(\infty)}} \right)^2$;
- 10 Compute $\tilde{a}_l = (h_0 \circ \dots \circ h_{k-1})(A_l)$ for $l = 0, \dots, m+2$, with branching $(-1)^{1/2} = i$;
- 11 Compute $\tilde{b}_l = (h_0 \circ \dots \circ h_{k-1})(B_l)$ for $l = 0, \dots, n+2$, with branching $(-1)^{1/2} = -i$;
- 12 Apply a Möbius transformation T that takes $\{\tilde{a}_{m+1}, \tilde{b}_{n+1}, \frac{1}{2}(\tilde{a}_{m+2} + \tilde{a}_{n+2})\}$ to $\{-1, 1, \infty\}$ on all the transformed points;

An illustration of the partial welding algorithm is given in Figure 7. As a remark, to weld two subdomains obtained by the local parameterization step partially, we only need to extract their boundary points on \mathbb{C} and apply Algorithm 2. The interior points of the two flattened subdomains are not needed. With the updated coordinates of the boundary points of the subdomains, we can then easily obtain the desired global conformal parameterization by solving a number of sparse linear systems. The details will be described in Section 4.5.

4.4. Enforcing additional constraints. Before moving on to the step of obtaining the final global parameterization, it is possible for us to include an optional step here and enforce additional constraints for achieving disk conformal parameterization and spherical conformal parameterization.

4.4.1. Constraints for disk conformal parameterization. If the input simply-connected surface \mathcal{S} is open, one can further restrict the target parameter domain to be the unit disk in the proposed method, thereby achieving a disk conformal parameterization. This is done by adding an extra step of applying the geodesic algorithm introduced in Section 3.4 to the global boundary $\partial\mathcal{S}$. Note that the points on $\partial\mathcal{S}$ are distributed into various subdomains. Therefore, we first extract the coordinates of those boundary points from the partial welding result. Once the mapping that takes those points to the unit circle is determined, we apply

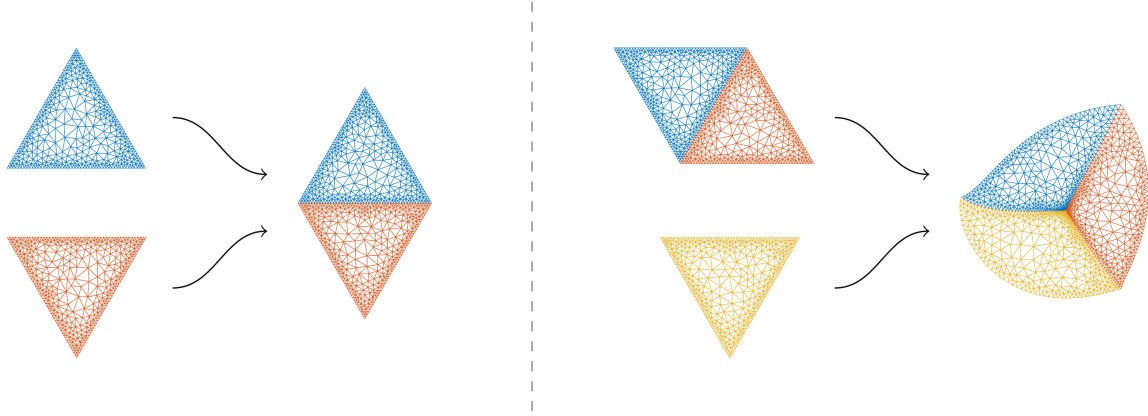


Figure 7. An illustration of partial welding. Suppose we are given a pyramid-like surface with three triangular faces (colored in blue, red, yellow). Each of them have been flattened onto the plane. In order to weld the boundaries of the three triangles, we apply Algorithm 2 twice. First, we glue the common boundaries of the blue and red triangles via partial welding. Then, we glue the common boundaries of the yellow triangle and the other two triangles via partial welding. Note that the interior points and the mesh structure of each triangle are plotted for a better visualization only. In the actual computation of partial welding, only the boundary points are involved.

the map for transforming the boundary coordinates of every flattened subdomain onto the unit disk. This results in boundary coordinates for the subdomains that yield a disk conformal parameterization upon solving the Laplace equation (details to be described in Section 4.5).

4.4.2. Constraints for spherical conformal parameterization. For genus-0 closed surfaces, one common choice of the parameter domain is the unit sphere \mathbb{S}^2 . In case the input surface \mathcal{S} is a genus-0 closed surface, we can modify our framework so that the partial welding procedure is repeated until two large components are left. Then, for the last welding, we use a closed welding instead of a partial welding to glue the entire boundaries of the two large components. As all boundaries are glued, the resulting boundary coordinates of the subdomains on the extended complex plane yield a spherical conformal parameterization upon solving the Laplace equation (details to be described in Section 4.5).

4.5. Obtaining the global conformal parameterization. After obtaining the new boundary constraints that satisfy the consistency condition, we can compute the global conformal parameterization of the input surface \mathcal{S} by finding a harmonic map $\tilde{\varphi}_i : \mathcal{S}_i \rightarrow \mathbb{R}^2$ for each submesh with the new boundary constraints. More explicitly, it suffices to solve the Laplace equation

$$(4.22) \quad \Delta \tilde{\varphi}_i = 0$$

subject to the new boundary constraints. Again, note that the computations for the K submeshes are independent and so this step is parallelizable. Because of the consistency between the boundaries of all subdomains, the new local parameterization results can be glued

seamlessly, thereby forming a global conformal parameterization. One can further ensure the bijectivity of each subdomain using the idea of quasi-conformal composition (see Section 3.5). More specifically, we compute the Beltrami coefficient of the inverse mapping $\tilde{\varphi}_i^{-1}$ (denoted by $\mu_{\tilde{\varphi}_i^{-1}}$). We can then determine whether $\tilde{\varphi}_i$ is folding-free by checking if $\|\mu_{\tilde{\varphi}_i^{-1}}\|_\infty > 1$ (or close to 1 in the discrete case). If so, we compose $\tilde{\varphi}_i$ with another mapping that is associated with the Beltrami coefficient $\mu_{\tilde{\varphi}_i^{-1}}$ to fix the fold-overs as guaranteed by quasi-conformal theory. With this additional step, we can ensure the bijectivity of the resulting global conformal parameterization.

Note that the resulting global parameterization lies in the extended complex plane. In case \mathcal{S} is a genus-0 closed surface, we add a stereographic projection step to convert it to a spherical parameterization. Algorithm 3 summarizes the proposed method.

Algorithm 3: Parallelizable global conformal parameterization of simply-connected surfaces (PGCP)

Input: A simply-connected surface mesh $\mathcal{S} = (\mathcal{V}, \mathcal{F})$, a set of edges $\tilde{\mathcal{E}}$ for the partition.

Output: A global conformal parameterization $\varphi : \mathcal{S} \rightarrow \mathbb{R}^2$ or \mathbb{S}^2 .

- 1 Partition the mesh into K submeshes based on $\tilde{\mathcal{E}}$;
 - 2 **for** $i = 1, \dots, K$ **do**
 - 3 Compute a conformal parameterization of $\mathcal{S}_i = (\mathcal{V}_i, \mathcal{F}_i)$ using DNCP. Only the boundary coordinates of the parameterization are kept;
 - 4 Perform partial welding as described in Algorithm 2 to update the boundary coordinates;
 - 5 (Optional) To achieve disk conformal parameterization, further apply the geodesic algorithm [40]. To achieve spherical conformal parameterization, perform conformal welding on the last two components obtained by partial welding;
 - 6 **for** $i = 1, \dots, K$ **do**
 - 7 Solve the Laplace equation $\Delta \tilde{\varphi}_i = 0$ with the new boundary constraints for each \mathcal{S}_i ;
 - 8 Compute the Beltrami coefficient $\mu_{\tilde{\varphi}_i^{-1}}$ to check whether $\tilde{\varphi}_i$ is folding-free. If not, fix the fold-overs in $\tilde{\varphi}_i$ using quasi-conformal composition;
 - 9 The solutions $\tilde{\varphi}_i$ for all \mathcal{S}_i together form a global conformal parameterization φ . For spherical conformal parameterization, further apply the stereographic projection to map the result onto \mathbb{S}^2 ;
-

As a remark, the novel combination of local parameterization and partial welding in our proposed method significantly improves the computational efficiency of global conformal parameterization. For a direct solver of global conformal parameterization, a computation in at least $O(|\mathcal{V}|^2)$ is needed (as a $2|\mathcal{V}| \times 2|\mathcal{V}|$ linear system is involved). By contrast, one can see that the interior parts of the submeshes are not used in the partial welding step in our method. Therefore, the heaviest computation involved in our proposed method only takes $O(|\mathcal{B}|^2)$, where \mathcal{B} is the collection of boundary points of the subdomains. The computation of the local parameterizations at the beginning and the harmonic maps at the end of our proposed method can both be parallelized, so that each computation runs in $O(|\mathcal{V}_i|^2)$.

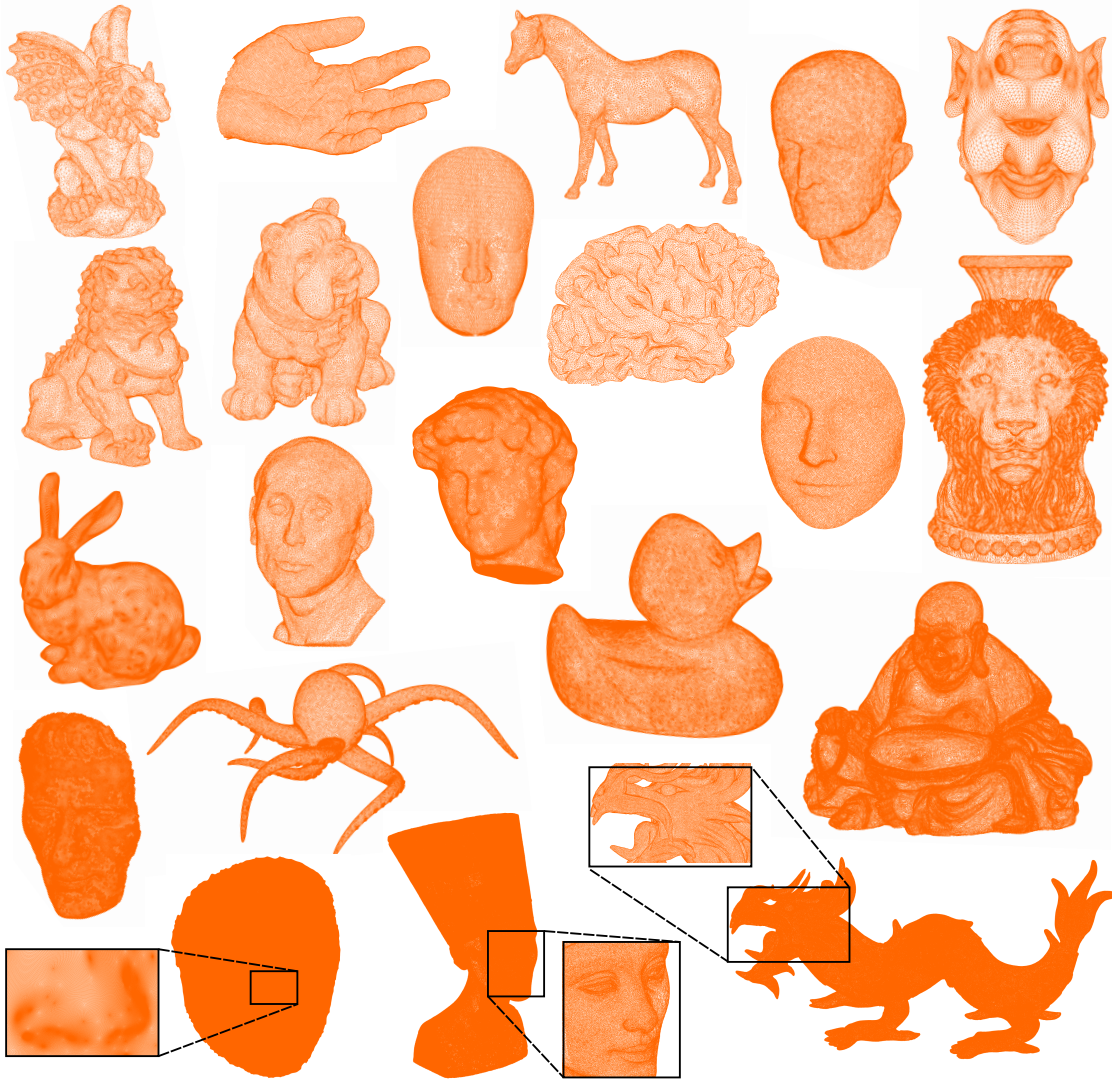


Figure 8. A gallery of simply-connected surface meshes used in our experiments. Our PGCP method is capable of handling a wide range of simply-connected surfaces with different geometry, mesh quality and resolution.

5. Experiments. Our proposed PGCP method is implemented in MATLAB, with the Parallel Computing Toolbox utilized for achieving parallelization. The sparse linear systems are solved using the backslash operator in MATLAB. All experiments are performed on a PC with Intel i7-6700K quad-core CPU and 16 GB RAM. To evaluate the performance of our proposed method, we adapt various simply-connected surface meshes from multiple free 3D model repositories [46–49] (see Figure 8). As for the distortion measure, we define the angular distortion of an angle $[v_i, v_j, v_k]$ (in degree) under the conformal parameterization φ by

$$(5.1) \quad d([v_i, v_j, v_k]) = \angle[\varphi(v_i), \varphi(v_j), \varphi(v_k)] - \angle[v_i, v_j, v_k].$$

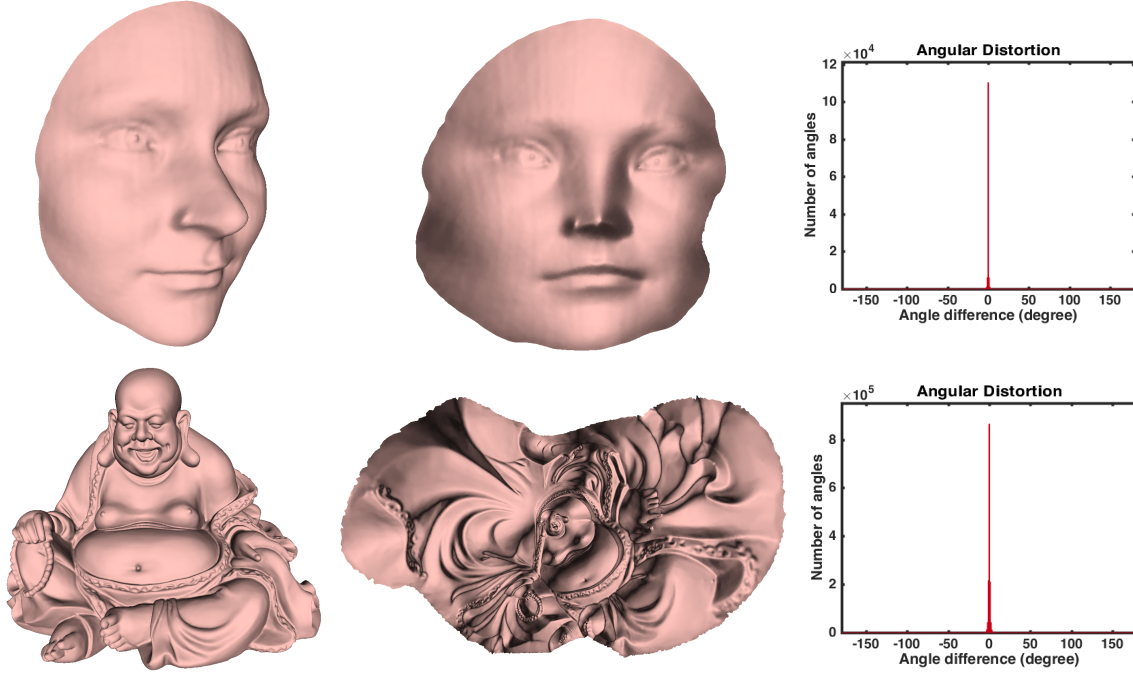


Figure 9. Free-boundary conformal parameterizations of simply-connected open surfaces obtained by our proposed PGCP method, rendered with normal map shader.

Surface	# vertices	SCP [17]		CETM [18]		PGCP	
		Time (s)	mean($ d $)	Time (s)	mean($ d $)	Time (s)	mean($ d $)
Sophie	21K	1.1	0.2	1.2	0.2	0.6	0.2
Niccolò da Uzzano	25K	1.3	0.6	Failed		0.7	0.6
Mask	32K	1.6	0.2	7.3	0.2	0.9	0.2
Max Planck	50K	2.6	0.5	5.5	0.5	1.5	0.5
Bunny	85K	4.4	0.5	18.8	0.5	2.0	0.5
Julius	220K	14.2	0.1	19.5	0.1	6.6	0.1
Buddha	240K	13.7	0.6	49.0	0.6	9.2	0.6
Face	1M	85.2	< 0.1	98.1	< 0.1	47.6	< 0.1

Table 1

The performance of spectral conformal parameterization (SCP) [17], conformal equivalence of triangle meshes (CETM) [18] and PGCP for free-boundary conformal parameterization of simply-connected open surfaces.

5.1. Free-boundary conformal parameterization of simply-connected open surfaces.

We first consider computing free-boundary global conformal parameterization of simply-connected open surfaces using our proposed PGCP method (see Figure 9 for examples). To assess the performance of our method, we compare it with the spectral conformal parameterization (SCP) [17] and conformal equivalence of triangle meshes (CETM) [18] in terms of the computation time and the angular distortion (see Table 1). The MATLAB version of SCP is

implemented by the authors, and the MATLAB version of CETM can be found at [50]. The experimental results show that our proposed method is significantly faster than both SCP and CETM by over 40% and 70% respectively on average, while maintaining comparable accuracy in terms of the average angular distortion. This demonstrates the effectiveness of our method for free-boundary global conformal parameterization.

5.2. Disk conformal parameterization of simply-connected open surfaces. Besides free-boundary global conformal parameterization, our proposed PGCP method can also achieve disk conformal parameterization of simply-connected open surfaces (see Figure 10 for examples). To evaluate the performance of our method, we compare it with the state-of-the-art linear disk conformal map (LDM) method [25] and the conformal energy minimization (CEM) method [26] (see Table 2). The MATLAB version of LDM can be found at [51], and the MATLAB version of CEM can be found at [52]. It can be observed that our method is significantly faster than LDM and CEM by over 50% and 30% on average respectively. Also, our method achieves comparable or smaller angular distortion when compared to the two other methods. This shows that our method is advantageous for disk conformal parameterization.

5.3. Spherical conformal parameterization of simply-connected closed surfaces. We then consider computing spherical conformal parameterization of genus-0 closed surfaces using our proposed PGCP method (see Figure 11 for examples). To evaluate the performance, we compare our proposed method with the state-of-the-art folding-free global conformal mapping (FFGCM) algorithm [30] and the FLASH algorithm [31] (see Table 3). The MATLAB version of FFGCM is kindly provided by the authors, and the MATLAB version of FLASH can be found at [53]. Because of the “divide-and-conquer” nature of our method, our method is capable of producing spherical conformal parameterizations with a smaller angular distortion when compared to the two state-of-the-art algorithms. In particular, the FLASH algorithm involves puncturing a triangle from the input surface and flattening the punctured surface onto a big triangular domain. This step unavoidably creates squeezed regions and produces certain angular distortions. While the distortions are alleviated in the subsequent step using quasi-conformal composition, the step again involves a domain where most vertices are squeezed at the interior, which leads to some distortions. By contrast, our proposed PGCP method flattens each submesh naturally, with the shape of the submesh boundary taken into consideration. This effectively reduces the angular distortions, thereby producing a spherical conformal parameterization with a better accuracy. Moreover, because of the ability of exploiting parallelism, our method achieves a significant reduction in computational time by over 90% on average when compared to FFGCM. When compared to FLASH, our method achieves comparable efficiency for moderate meshes and a notable reduction in computational time by around 25% for dense meshes. This shows the advantages of our method for spherical conformal parameterization.

5.4. Applications. The above experiments demonstrate the improvement of our proposed PGCP method over the state-of-the-art conformal parameterization algorithms. In this section, we discuss the applications of it.

5.4.1. Texture mapping. One application of our proposed PGCP method is texture mapping. After conformally flattening a surface onto the plane using our method, we can

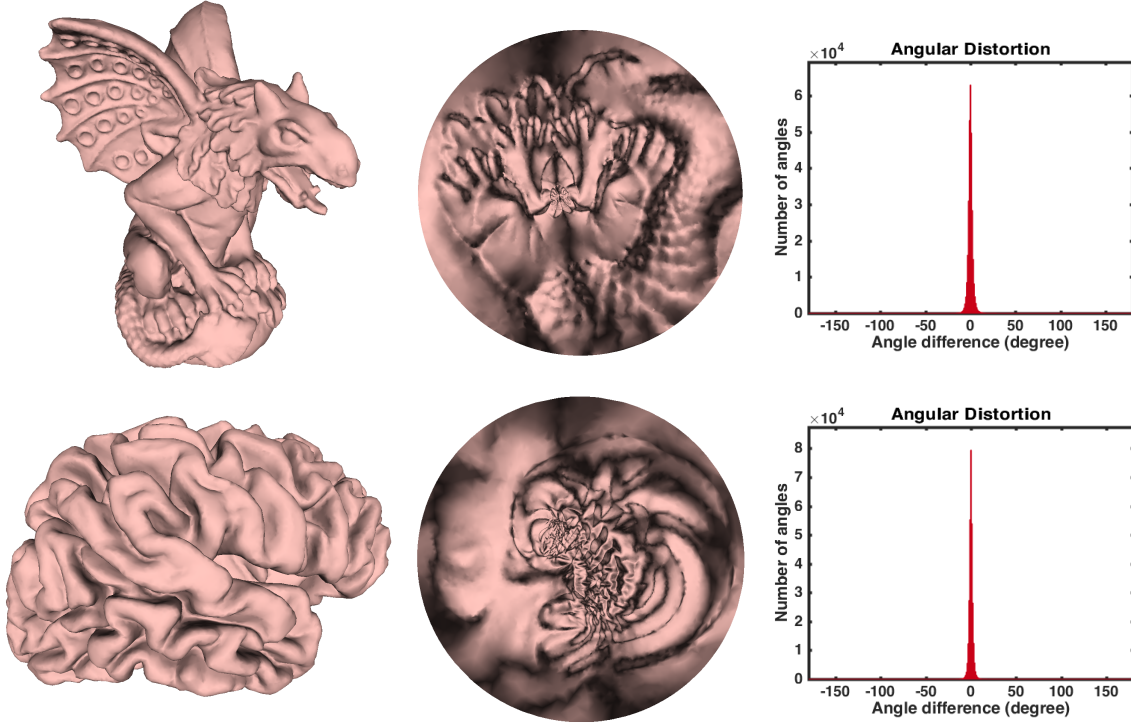


Figure 10. Disk conformal parameterizations of simply-connected open surfaces obtained by our proposed PGCP method, rendered with normal map shader.

Surface	# vertices	LDM [25]		CEM [26]		PGCP	
		Time (s)	mean($ d $)	Time (s)	mean($ d $)	Time (s)	mean($ d $)
Ogre	20K	1.1	1.5	0.3	2.6	0.5	1.5
Niccolò da Uzzano	25K	1.6	0.8	1.4	1.3	0.8	0.8
Brain	48K	2.9	1.6	2.9	1.5	1.3	1.5
Gargoyle	50K	3.1	1.9	2.8	2.1	1.4	1.9
Hand	53K	3.4	1.2	3.4	1.2	1.4	1.2
Octopus	150K	15.4	7.2	10.4	24.0	8.9	5.6
Buddha	240K	22.4	0.7	25.1	0.7	11.4	0.7
Nefertiti	1M	87.9	2.9	83.2	4.2	52.7	2.9

Table 2

The performance of linear disk conformal map (LDM) [25], conformal energy minimization (CEM) [26] and PGCP for disk conformal parameterization of simply-connected open surfaces.

design a texture on the parameter domain. Since there is a 1-1 correspondence between the input surface and the parameter domain, we can then use the inverse mapping to map the texture back onto the surface, thereby obtaining a surface with the desired texture on it. Several examples are shown in Figure 12. It is noteworthy that our method is conformal

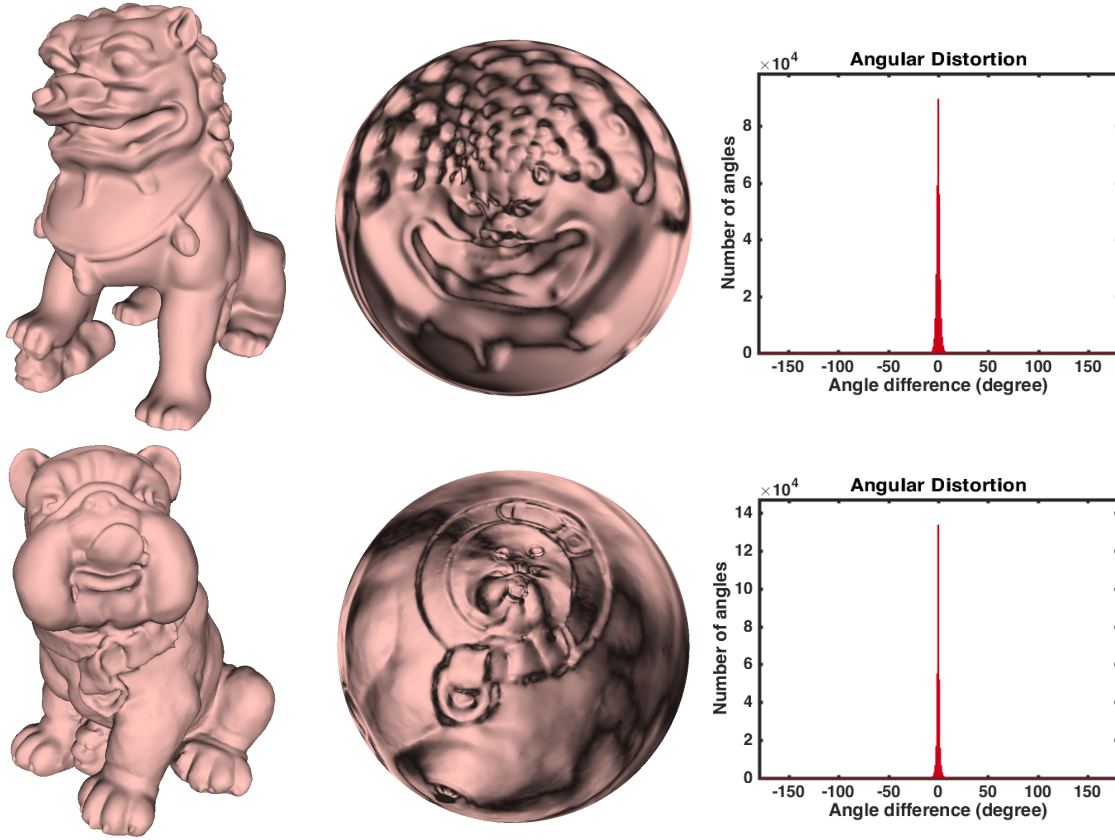


Figure 11. Spherical conformal parameterizations of genus-0 closed surfaces obtained by our proposed PGCP method, rendered with normal map shader.

Surface	# vertices	FFGCM [30]		FLASH [31]		PGCP	
		Time (s)	mean(d)	Time (s)	mean(d)	Time (s)	mean(d)
Horse	20K	12.1	11.0	0.4	3.0	0.4	2.7
Bulldog	50K	22.0	1.0	0.9	1.1	1.0	1.0
Chinese Lion	50K	29.3	1.3	1.1	1.3	1.1	1.3
Duck	100K	100.4	1.1	2.2	0.4	2.4	0.3
David	130K	46.6	0.2	3.5	0.2	3.4	0.2
Octopus	150K	112.3	37.2	10.1	6.9	7.1	2.6
Lion Vase	210K	222.7	14.4	4.5	0.8	4.7	0.7
Asian Dragon	1M	Failed		64.4	1.3	48.5	0.9

Table 3

The performance of folding-free global conformal mapping (FFGCM) [30], FLASH [31] and PGCP for spherical conformal parameterization of genus-0 closed surfaces.

and hence the local geometry of the texture pattern is well preserved. For instance, the checkerboard texture shown in Figure 12 can maintain its orthogonality on the Ogre surface.

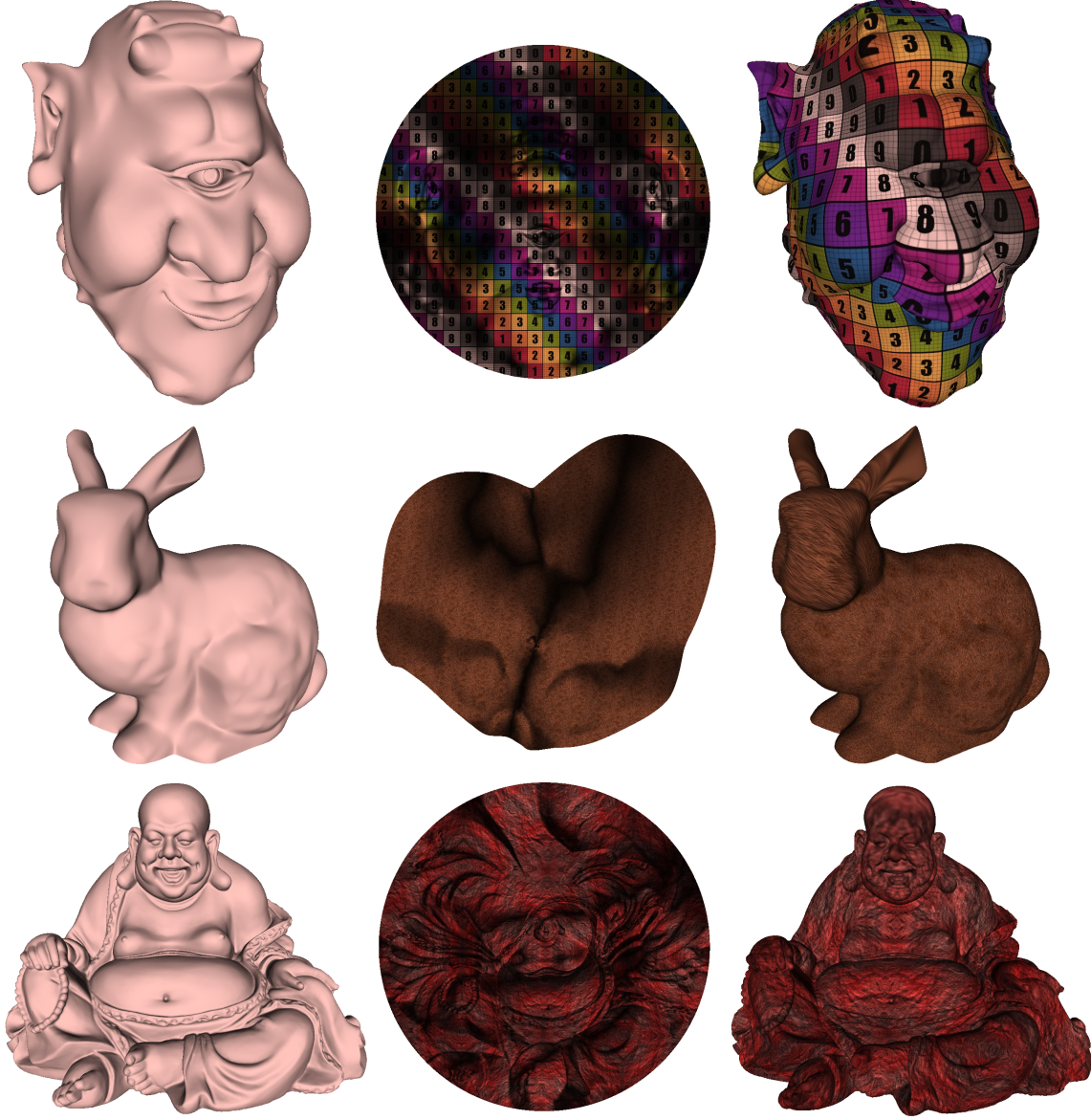


Figure 12. Texture mapping via our proposed PGCP method. Left: the input mesh. Middle: the parameterization achieved by PGCP rendered with normal map shader, overlaid with a texture (colored checkerboard/hair/stone). Right: the texture mapping result.

5.4.2. Surface remeshing. Our proposed PGCP method can also be applied to surface remeshing, which aims at improving the mesh quality of a given surface. By conformally parameterizing the surface and constructing a regular mesh structure on the parameter domain, we can use the inverse mapping to map the mesh structure back onto the surface,

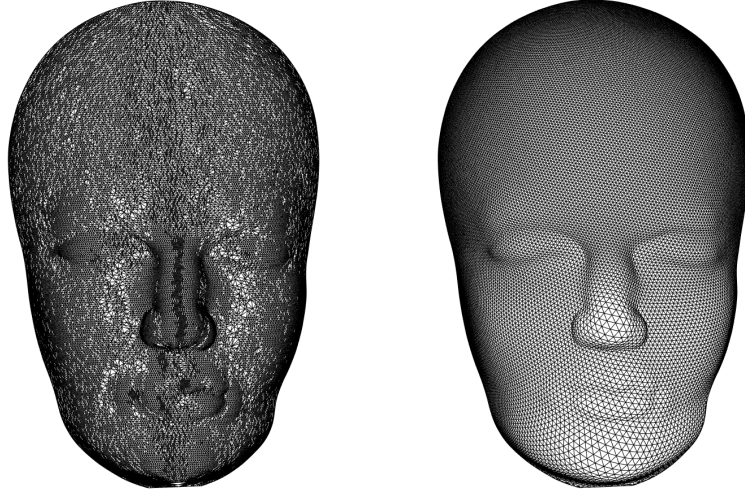


Figure 13. Surface remeshing via our proposed PGCP method. Left: the input surface. Right: the remeshed surface with improved mesh quality.

thereby remeshing the surface (see Figure 13 for example). It is noteworthy that since the parameterization is conformal, the regularity of the mesh structure defined on the parameter domain is well-preserved on the surface.

5.4.3. Solving PDEs on surfaces. Another notable application of our proposed PGCP method is solving PDEs on surfaces [42]. While solving PDEs on a general surface is difficult, solving them on a standard parameter domain such as the unit sphere or the unit disk is relatively easy. Figure 14 shows an example of patterns formed on the genus-0 David surface by solving the time-dependent Ginzburg-Landau equation on the spherical conformal parameterization obtained by our proposed PGCP method. The PDE on the sphere is solved using Chebfun [54]. The example demonstrates the use of our method for PDE-based surface decoration.

5.4.4. Other applications. Some other possible applications of conformal parameterizations include surface registration [31], medical visualization [24] and surface morphing [26]. As our proposed PGCP method is advantageous over the state-of-the-art algorithms in both the computational time and the conformal distortion, these tasks can be done with higher efficiency and accuracy using our method.

6. Conclusion. In this work, we have proposed a novel parallelizable global conformal parameterization method called PGCP for simply-connected surfaces. Given a triangle mesh, we partition it into submeshes and conformally flatten each of them using DNCP. As the local parameterization results do not yield a consistent global parameterization, we extract their boundary points to integrate them using a novel technique called partial welding. Using the modified boundaries for all submeshes, harmonic maps can be computed to yield a global conformal parameterization, with bijectivity guaranteed by quasi-conformal theory. Additional steps can be included to produce disk conformal parameterizations for surfaces with boundary,

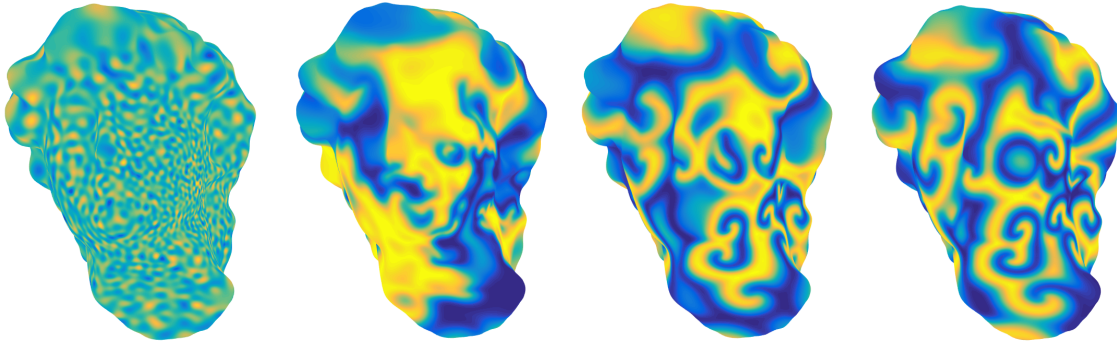


Figure 14. Patterns formed on the genus-0 David surface by solving the Ginzburg-Landau equation on the spherical conformal parameterization. The leftmost is the initialization, and the rightmost is the final result.

and spherical conformal parameterizations for genus-0 closed surfaces.

Most parts of our proposed method, such as the initial local conformal parameterization step and the last harmonic mapping step, can be computed independently in a distributed manner. The only global computation involved in our algorithm takes merely boundary data points of the submeshes, which are much fewer than the vertices of the entire mesh. Experimental results have demonstrated the significant improvement in efficiency and accuracy achieved by our proposed method when compared to the state-of-the-art approaches for free-boundary conformal parameterization, disk conformal parameterization and spherical conformal parameterization.

For future work, we plan to explore the possibility of extending our method for quasi-conformal parameterizations and mappings [43–45]. More specifically, note that the partial welding step in our proposed method is conformal, and the quasi-conformal dilatation of a map is preserved under the composition with conformal maps. Therefore, it should be possible for us to compute quasi-conformal parameterizations and mappings for dense meshes by a combination of local quasi-conformal maps of submeshes and partial welding. Another possible future work is the extension of our method for point clouds. As the partial welding approach uses only the boundary data points of the flattened submeshes but not the mesh structure of them, it should also be applicable for subdomains of a point cloud. Combining the partial welding approach with some existing conformal parameterization methods for disk-type point clouds will then yield a parallelizable global conformal parameterization method for point clouds.

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