# Non-self-overlapping structured grid generation on an $n$-sided surface 

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#### Abstract

Most existing meshing algorithms for a 2D or shell figure requires the figure to have exactly 4 sides. Generating structured grids in the $n$-sided parametric region of a trimmed surface thus usually requires to first partition the region into 4 -sided sub-regions. We address the automatic structured grid generation problem in an $n$-sided region by fitting a planar Gregory patch so that the partition requirement is naturally avoided. However, selfoverlapping may occur in some portions of the algebraically generated grid; this severely limits its usage in most of engineering and scientific applications where a grid system with no self-intersecting is strictly required. To solve the problem, we use a functional optimization approach to move grid nodes in the $u-v$ domain of the trimmed surface to eliminate the self-overlapping. The derivatives of a Gregory patch, which are extremely difficult to compute analytically, are not required in our method. Thus, our optimization algorithm compares favorably at least in terms of speed with some other mesh optimization algorithms, such as elliptic PDE method. In addition, to overcome the difficulty of guessing a good initial position of every grid node for the conjugate gradient method, a progressive optimization algorithm is incorporated in our optimization. Experiment results are given to illustrate the usefulness and effectiveness of the presented method.


Keywords: self-overlapping, structured grid, quadrilateral grid, Gregory patch, $n$-sided, and trimmed surface.

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## 1. Introduction

In computer-aided engineering, geometric modeling, computer graphics, and many other applications, trimmed parametric surfaces are widely adopted [1]. After a parametric surface $S$ intersects with other surfaces, only a portion of the surface patch is used in defining a meaningful shape, which is called a trimmed (parametric) surface $S_{T}$. $S_{T}$ is constrained by the same mathematical surface equation as $S(u, v)$, but its parametric domain is only a portion of that of $S$. The parametric area of $S_{T}$, to be denoted as $\mathrm{P}_{S_{T}}$, lies inside $(u, v) \in[0,1] \times[0,1]$ (assuming the $u-v$ domain of $S$ is normalized) and is bounded by a number of curves (see Fig. 1-1). Each boundary curve of $\mathrm{P}_{S_{T}}$ is expressed as a parametric equation of the form $b_{i}=\left[\begin{array}{ll}u_{i}(t) & v_{i}(t)\end{array}\right]$, where $t \in[0,1]$. Formally, a trimmed $n$-sided surface patch is defined below.

Definition 1-1 A trimmed parametric surface $S_{T}$, whose $n$ boundary curves form a Jordan curve in its parametric area $\mathrm{P}_{S_{T}}$, is defined as a trimmed n-sided (parametric) surface patch.

(a) an example surface

(b) the parametric area in $u-v$ domain

Fig. 1-1 A trimmed $\boldsymbol{n}$-sided surface and its parametric area

An example of a trimmed $n$-sided patch and its parametric area is given in Fig. 1-1. The task of approximating a trimmed surface by a complex of simple planar elements (either triangular or quadrilateral) plays an important role in engineering computing; this is referred to as the surface meshing operation, which has been studied for many years [2-16]. The two most powerful analysis tools in engineering are the finite element methods and the finite difference methods. The finite element method usually adopts either triangular grids or quadrilateral grids, and the grids can be structured or unstructured. However, the finite difference method generally uses structured quadrilateral grids (or simply called structured grids). The structured grids can be generated algebraically or as the solution of Partial Deferential Equations (PDEs). Surface grid generation algorithms are based on a projection strategy, consisting in generating a grid on the surface parametric plane, and then to project it on the surface. However during the projection step the grid characteristics, such as point distribution and orthogonality, are often not correctly transported. In this approach, we solve the structured grid generation problem in an $n$-sided region while preserving the non-overlapping property in a numerical optimization manner.

Within all the surface grid generation approaches, algebraic grid generation is some form of interpolation from boundary points - different approaches use different kinds of interpolation [3-5]. Overlapping may, however, happen on some portions of algebraically generated grids, which must be corrected in order for the mesh to be useable for almost any application. Also, error may be generated when converting generic boundary curves into curves with a specified representation. Grid generation is actually a boundary-value problem, so grids can be generated from point distribution on boundaries by solving elliptic PDEs in the field [6-10]. The smoothness properties and extremum principles of some such PDE systems can serve to produce smooth grids without boundary overlapping. The most recent work of Arina [11] overcame the difficulties of point distribution and orthogonality on the mapped a surface by a conformal map, preserving angles and scale-length ratios. His method consists of two major steps: the parameterization of surface by isothermal coordinates; the generation of a two-dimensional grid on the conformal parametric plane and its projection on the surface. All the above approaches however cannot be directly applied to a region with $n$ sides (when $n>4$ ), so the $n$-sided region has to be partitioned into several non-overlapping blocks with 4 sides. For the region with a complex shape (as shown in Fig. 1-2a and 1-2b), generating such partitions is not a straightforward work - skeleton structure $[12,13]$ or background triangulation $[14,15]$ may be required to construct. However, their resultant grids are usually unstructured.

In this paper, a novel method is presented for generating structured quadrilateral grids with no selfoverlapping on a trimmed $n$-sided patch by automatically fitting non-self-overlapping multi-block planar grids into the parametric region of the given surface. The idea is to first construct an initial structured algebraic grid based on the Gregory patch mapping [16] and then eliminate any possible self-overlapping on the mesh by performing a functional optimization. In our approach, a planar Gregory patch $G(\xi, \eta)$ is adopted to fit an algebraic grid in the $u-v$ domain of $S_{T}$; and since the $\xi-\eta$ domain $\mathrm{P}_{G}$ of $G(\xi, \eta)$ is a regular $n$-sided polygon, it is easy to divide $\mathrm{P}_{G}$ into $n 4$-sided sub-regions and grid each sub-region automatically (see Fig. 12 c and $1-2 \mathrm{~d})$. The mapping between the $\xi-\eta$ domain and the $u-v$ domain by a Gregory patch nevertheless may generate overlapping (see Fig. 1-2e and 1-2f). To mend that, we develop a functional optimization method to eliminate such overlapping in the $u-v$ domain; the shapes of grid elements in the $u-v$ space are also adjusted in the optimization. One question arising naturally is: why not using PDEs for this optimization? Our answer is that, when enhancing the quality of grids by the PDE methods, the derivatives such as $G_{\xi}, G_{\eta}, G_{\xi \xi}$, $G_{\eta \eta}$, and $G_{\xi \eta}$ are needed; for a Gregory patch (detail in section 3), it is hard to give the analytical formulas for these terms, and computing them numerically is a very time-consuming process. Our method compares favorably to PDEs in this regard, as only the positions of every grid nodes in the $u-v$ domain are required. Thus, the speed of our method is faster.

The paper is organized as follows. In section 3, after giving some necessary definitions and preliminaries for the Gregory patch mapping, we introduce the method to determine the initial algebraic grid by a planar Gregory patch. We then formulate the self-overlapping problem in the $u-v$ domain as a singularity problem in the derivatives on the Gregory patch, in section 4.1. Based on this formulation, the corresponding objective function of our optimization is then derived in section 4.2, with a shape control term also added to the objective function. The details of the numerical implementation of the optimization are given in section 4.3. As our numerical implementation is iterative, to overcome the difficulty of "guessing" a good initial grid, the idea of
progressive optimization is introduced in section 4.4. Finally, some experimental results are shown in section 5 to demonstrate the power of our approach, with some conclusion remarks offered in the last section.

(a) a trimmed $\boldsymbol{n}$-sided surface $S_{T}(u, v)$ in the

$$
x-y-z \text { space }
$$


(c) the parametric area of $G(\xi, \eta)$ in the $\xi-\eta$ domain

(e) the grid in the $u-v$ domain (with overlapping)

(b) the parametric area of $S_{T}(u, v)$ in the $u-v$ domain

(d) the grid in the $\xi-\eta$ domain

(f) the grid for the surface $S_{T}$ (with overlapping)

Fig. 1-2 Mapping from the $x-y-z$ space to the $\xi-\eta$ domain via the $u-v$ domain (Example I)

## 2. Gregory patch mapping

The necessary definitions and preliminaries of a Gregory patch are first given here.


Fig. 2-1 Define a Gregory corner interpolator
Definition 2-1 Let $P(u): 0 \leq u \leq 1$ and $Q(v): 0 \leq v \leq 1$ be two regular curves in $\mathfrak{R}^{3}$ with $P(0)=Q(0)$, and $T_{P}(u): 0 \leq u \leq 1$ and $T_{Q}(v): 0 \leq v \leq 1$ be two $C^{1}$ vector functions in $\Re^{3}$ satisfying $T_{P}(0)=\left.\frac{d Q(v)}{d v}\right|_{v=0}$ and $T_{Q}(0)=\left.\frac{d P(u)}{d u}\right|_{u=0}$, the Gregory corner interpolator of the four, $\left\{P(u), Q(v), T_{P}(u), T_{Q}(v)\right\}$, is a surface in $\mathfrak{R}^{3}$ defined by

$$
\begin{equation*}
r(u, v)=P(u)+v T_{P}(u)+Q(v)+u T_{Q}(v)-P(0)-v T_{P}(0)-u T_{Q}(0)-u v \frac{v T_{P}^{\prime}(0)+u T_{Q}^{\prime}(0)}{u+v} \tag{1}
\end{equation*}
$$

The Gregory corner interpolator function $r(u, v)$ agrees with $P(u)$ and $Q(v)$ along the two sides (i.e., $r(u, 0)=P(u)$ and $r(0, v)=Q(v))$. Also, its partial derivatives with respect to $u$ and $v$ agree with $T_{P}(u)$ and $T_{Q}(v)$ along the respective sides $-\left.\frac{\partial r(u, v)}{\partial v}\right|_{v=0}=T_{P}(u)$ and $\left.\frac{\partial r(u, v)}{\partial u}\right|_{u=0}=T_{Q}(v)$ since $T_{P}(0)=Q^{\prime}(0)$ and $T_{Q}(0)=P^{\prime}(0)$. For an $n$-sided 3D surface, $n$ such interpolator functions can be defined on the $n$ corners; the final surface is the weighted sum of the $n$ functions [17, 18]. The details are defined as follows.

Definition 2-2 The parametric domain of a Gregory patch with $n$ sides is defined as a unit length regular $n$ gon in the $\xi-\eta$ domain.

We name the parametric domain of a Gregory patch $G$ as $\mathrm{P}_{G}$, where all corners $X_{k}(k=0,1, \ldots, n-1)$ are ordered in the anti-clockwise (as shown in Fig. 2-2). Given a point $X=\left(\xi_{0}, \eta_{0}\right)$ inside $\mathrm{P}_{G}$, when computing its three dimensional position defined by a Gregory corner interpolator $r_{k}\left(u_{k}, v_{k}\right)$, the parameters $\left(u_{k}, v_{k}\right)$ of the point corresponding to the $k$ th corner $X_{k}$ are defined as

$$
\begin{equation*}
\left(u_{k}, v_{k}\right)=\left(\frac{d_{k-1}}{d_{k-1}+d_{k+1}}, \frac{d_{k}}{d_{k-2}+d_{k}}\right) \tag{2}
\end{equation*}
$$

where $d_{k}$ represents the perpendicular distance from $X$ to the side $X_{k} X_{k+1}$. It is easy to find that if $\left(\xi_{0}, \eta_{0}\right)$ lies on the side $X_{k} X_{k+1}, v_{k}=0$ since $d_{k}=0$; if $\left(\xi_{0}, \eta_{0}\right)$ is on $X_{k-1} X_{k}, u_{k}=0$ since $d_{k-1}=0$; when $\left(\xi_{0}, \eta_{0}\right)$ and $X_{k+1}$ coincides, we have $u_{k}=1$ by equation (2); and when $\left(\xi_{0}, \eta_{0}\right)$ and $X_{k-1}$ coincides, we have $v_{k}=1$.


Fig. 2-2 $\mathrm{P}_{G}$ of a Gregory patch with five sides


Fig. 2-3 Define a Gregory patch

Definition 2-3 If $C_{0}(u), C_{1}(u), \ldots, C_{n-1}(u)$ are $n$ regular 3D curves that form a closed loop in three dimensional space, that is $C_{k}(1)=C_{(k+1) \bmod n}(0)(k=0,1, \ldots, n-1)$, and $T_{C_{0}}(u), T_{C_{1}}(u), \ldots, T_{C_{n-1}}(u)$ are $n$ continuous 3 D vector functions defined on the $C_{k}(u) \mathrm{s}$ respectively, the Gregory patch of $C_{k}(u) \mathrm{s}$ and $T_{C_{k}}(u) \mathrm{s}$ is defined as a mapping from $\mathrm{P}_{G}$ to $\mathfrak{R}^{3}$

$$
\begin{equation*}
G(X)=\sum_{k=0}^{n-1} w_{k}(X) r_{k}\left(u_{k}(X), v_{k}(X)\right) \tag{3}
\end{equation*}
$$

where $w_{k}(X)=\frac{\prod_{j \neq k-1, k} d_{j}^{2}}{\sum_{l=0}^{n-1} \prod_{j \neq l-1, l} d_{j}^{2}}$, and $r_{k}\left(u_{k}, v_{k}\right)$ represents the Gregory corner interpolator function for the $k$ th
corner of the four items $\left\{C_{k}(u), \bar{C}_{k}(v), T_{C_{k}}(u), \bar{T}_{C_{k}}(v)\right\}, \bar{C}_{k}(u)=C_{k}(1-u)$, and $\bar{T}_{C_{k}}(u)=T_{C_{k}}(1-u)$

Note that $w_{k}(X)$ is unity at the vertex $X_{k}$ and is zero on those edges of the $n$-gon not incident with $X_{k}$. Thus, it can be verified that the Gregory patch $G(X)$ is boundary conforming. That is, if an $X$ is on the boundary of $\mathrm{P}_{G}$, then $G(X)$ must be on one of the $C_{k}(u) \mathrm{s}$, and conversely, for any point $p \in \mathfrak{R}^{3}$ on any $C_{k}(u)$, there must be an $X$ on the boundary of $\mathrm{P}_{G}$ such that $p=G(X)$. If the $n$ boundary curves of a Gregory patch $G(X)$ all lie in a common plane, obviously $G(X)$ also lies in that plane, i.e., it is a planar Gregory patch. Next, we elaborate on how the Gregory patch mapping can be utilized to mesh an $n$-side trimmed parametric surface patch.

## 3. Algebraic grid generation based on Gregory mapping

Similar to other algebraic grid generation methods, the algebraic grid construction in our approach also consists of three steps: 1) forward mapping; 2) grid generation; and 3) inverse mapping. The forward mapping is
the mapping of the three dimensional physical surface $S_{T}$ to its underlying parametric area $\mathrm{P}_{S_{T}}$. By fitting a planar Gregory patch $G$ into $\mathrm{P}_{S_{T}}, \mathrm{P}_{S_{T}}$ of the surface $S_{T}$ is further mapped into the parametric $n$-gon $\mathrm{P}_{G}$ of $G$. In this planar Gregory patch $G$, all the $C_{k}(u) \mathrm{s}$ and $T_{C_{k}}(u) \mathrm{s}$ are determined by the 2 D boundary curves of $\mathrm{P}_{S_{T}}$ in the parametric domain instead of the 3D boundary curves of $S_{T}$. Grids will be generated in $\mathrm{P}_{G}$ and then mapped back into $\mathfrak{R}^{3}$, generating a structured grid system of $S_{T}$.

In the $\xi-\eta$ domain of $G$, the coordinates of parametric $n$-gon $\mathrm{P}_{G}$ 's corners are defined by

$$
\begin{equation*}
X_{k}=\left(\cos \frac{2 k \pi}{n}, \sin \frac{2 k \pi}{n}\right), \tag{4}
\end{equation*}
$$

where $n$ is the number of corners, and $k=0, \ldots, n-1$; so the position of middle points on the $n$ boundary sides can be computed by

$$
\begin{equation*}
E_{k}=\left(\frac{1}{2}\left(\cos \frac{2 k \pi}{n}+\cos \frac{2(k+1) \pi}{n}\right), \frac{1}{2}\left(\sin \frac{2 k \pi}{n}+\sin \frac{2(k+1) \pi}{n}\right)\right) . \tag{5}
\end{equation*}
$$

By the $X_{k} \mathrm{~s}, E_{k} \mathrm{~s}$, and the origin $o$, the $n$-gon $\mathrm{P}_{G}$ is divided into $n$ blocks. When establishing the $M \times N$ grid in the $k$ th sub-domain of $\mathrm{P}_{G}$, the coordinate of every grid node ( $\xi_{i, j}^{k}, \eta_{i, j}^{k}$ ) is determined by (as illustrated in Fig. 3-1)

$$
\left\{\begin{array}{l}
\xi_{i, j}^{k}=\left(1-\frac{i}{M}\right)\left[\left(1-\frac{j}{N}\right) \xi\left(X_{K}\right)+\frac{j}{N} \xi\left(E_{K-1}\right)\right]+\frac{i}{M}\left[\left(1-\frac{j}{N}\right) \xi\left(E_{K}\right)+\frac{j}{N} \xi(o)\right]  \tag{6}\\
\eta_{i, j}^{k}=\left(1-\frac{i}{M}\right)\left[\left(1-\frac{j}{N}\right) \eta\left(X_{K}\right)+\frac{j}{N} \eta\left(E_{K-1}\right)\right]+\frac{i}{M}\left[\left(1-\frac{j}{N}\right) \eta\left(E_{K}\right)+\frac{j}{N} \eta(o)\right]
\end{array},\right.
$$

where $\xi(\cdots)$ and $\eta(\cdots)$ represent the $\xi$ and $\eta$ coordinates of a point in the $\xi-\eta$ plane. In our approach, the $n$ sub-domains are meshed with the same number of $M$ and $N$, so the boundary nodes of adjacent sub-domains are coincident. To benefit the later grid optimization algorithm, we store the topological structure of the final grid $M^{G}$ by a pair of complex $\left(V^{G}, K^{G}\right)$, where $V^{G}$ is the set of grid nodes (the coincident vertices can be stored only once in $V^{G}$ ), and $K^{G}$ is a simplex complex specifying the connectivity of the grid simplices (the vertex-face and vertex-vertex adjacency information). Therefore, every node lying on the boundary of $\mathrm{P}_{G}$ has two incident faces, and the node coinciding with $o$ has $n$ incident faces, while all other inner nodes have four incident faces.


Fig. 3-1 Multi-block grids generation in the $\xi-\eta$ domain in $\mathrm{P}_{G}$

After the grid $M^{G}$ is constructed in $\mathrm{P}_{G}$, the coordinates of grid nodes are mapped backwards into $\mathfrak{R}^{3}$ by the algebraic equation of $G$ and $S_{T}$. However, when fitting a planar Gregory patch $G$ into the parametric domain $\mathrm{P}_{S_{T}}$ of $S_{T}$ in the $u-v$ plane, only the boundary curves - $C_{k}(u) \mathrm{s}$ are given by $\mathrm{P}_{S_{T}}$. To determine the algebraic equation of $G$ (Eq. (3)), the $T_{C_{k}}(u) \mathrm{s}$ are also needed. Here, a linear blending function is chosen. By the compatibility conditions at corners given in definition 2-1, we have

$$
T_{C_{i}}(0)=\left.\frac{d \bar{C}_{i-1}(u)}{d u}\right|_{u=0}=-\left.\frac{d C_{i-1}(u)}{d u}\right|_{u=1} \text { and } T_{C_{i}}(1)=\left.\frac{d C_{i+1}(u)}{d u}\right|_{u=0},
$$

the function of $T_{C_{i}}(u)$ is then given as

$$
\begin{equation*}
T_{C_{i}}(u)=(1-u) T_{C_{i}}(0)+u T_{C_{i}}(1) . \tag{7}
\end{equation*}
$$

Now, the final coordinates of grid nodes on the physical trimmed space can be expressed in the following form

$$
S_{T_{m}}=S\left(u\left(G\left(\xi_{m}, \eta_{m}\right)\right), v\left(G\left(\xi_{m}, \eta_{m}\right)\right)\right)
$$

where the functions $u(\cdots)$ and $v(\cdots)$ represent the $u$ and $v$ coordinates of a point on the Gregory patch.
As already exemplified in Fig. 1-2, however, when the boundary of $\mathrm{P}_{S_{T}}$ is complicated and convoluted, self-overlapping may occur in the $u-v$ plane by the planar Gregory patch mapping, which leads to a selfintersecting grid on $S_{T}$. In the following section, we introduce an iterative functional minimization method for the purpose of eliminating or reducing the self-overlapping in the $u-v$ domain.

## 4. Grid optimization

Once an initial structured mesh $M^{G}$ is generated on $\mathrm{P}_{S_{T}}$ by means of Gregory patch mapping, we next proceed to eliminate any possible self-overlapping in the mesh. This is achieved by first modeling this elimination process as a functional minimization problem and then introducing the formulas to optimize the derived objective function. To avoid "guessing" a good initial value in the iterative optimization process, a progressive optimization algorithm is also proposed.

### 4.1 Non-self-overlapping property

First, it is necessary to give a formal mathematical characterization of self-overlapping. Let us add a virtual axis $w$ perpendicular to the $u-v$ plane as $w=u \times v$. A Gregory patch $G(\xi, \eta)$ now is a planar region embedded in the $u \times v \times w$ space, denoted as $G(\xi, \eta)=[U(\xi, \eta) \quad V(\xi, \eta) \quad W(\xi, \eta)]^{T}$, where $U(\xi, \eta), V(\xi, \eta)$, and $W(\xi, \eta)$ are the components of $G(\xi, \eta)$ on the $u, v$, and $w$ axis respectively (note that $W(\xi, \eta)$ is a constant zero). We define the unit "normal vector" at any point $\left(\xi_{0}, \eta_{0}\right)$ on the patch $G(\xi, \eta)$ as $N\left(\xi_{0}, \eta_{0}\right)=\left.\frac{G_{\xi} \times G_{\eta}}{\left\|G_{\xi} \times G_{\eta}\right\|}\right|_{\left(\xi_{0}, \eta_{0}\right)}$. A point $\left(\xi_{0}, \eta_{0}\right)$ is said to be singular if its corresponding $\left\|G_{\xi} \times G_{\eta}\right\|$ is a zero vector. The lemma below is important as it stipulates the condition for guaranteeing the non-self-overlapping property.

Lemma 4-1 The mapping $G(\xi, \eta)=\left[\begin{array}{lll}U(\xi, \eta) & V(\xi, \eta) \quad W(\xi, \eta)\end{array}\right]^{T}$ has no self-overlapping if and only if there is no any singular point in the $n$-gon $\mathrm{P}_{G}$.

Proof. Let us argument the $G(\xi, \eta)$ by $G^{*}(\xi, \eta)=\left[U^{*}(\xi, \eta) \quad V^{*}(\xi, \eta) \quad W^{*}(\xi, \eta)\right]^{T}$ with $U^{*}(\xi, \eta)=U(\xi, \eta)$, $V^{*}(\xi, \eta)=V(\xi, \eta)$, and $W^{*}(\xi, \eta)=a \xi+b \eta$ for some real number $a$ and $b$. By properly choosing $a$ and $b$, one can enforce $G^{*}(\xi, \eta)$ to have non-zero length normal vector everywhere, thus it is a smooth regular surface in the $u \times v \times w$ space. Suppose first that $G(\xi, \eta)$ is self-overlapped. This means that there exist two distinct pairs $\left(\xi_{0}, \eta_{0}\right) \in \mathrm{P}_{G}$ and $\left(\xi_{1}, \eta_{1}\right) \in \mathrm{P}_{G}$, such that $\left(U^{*}\left(\xi_{0}, \eta_{0}\right), V^{*}\left(\xi_{0}, \eta_{0}\right)\right)=\left(U^{*}\left(\xi_{1}, \eta_{1}\right), V^{*}\left(\xi_{1}, \eta_{1}\right)\right)$. By properly selecting $a$ and $b$, one can also ensure that $W^{*}\left(\xi_{0}, \eta_{0}\right) \neq W^{*}\left(\xi_{1}, \eta_{1}\right)$. Let us intersect $G^{*}(\xi, \eta)$ with a plane $\Pi$ that is parallel to the $u-w$ plane and contains the two points

$$
p_{0}=\left(U^{*}\left(\xi_{0}, \eta_{0}\right), V^{*}\left(\xi_{0}, \eta_{0}\right), W^{*}\left(\xi_{0}, \eta_{0}\right)\right) \text { and } p_{1}=\left(U^{*}\left(\xi_{1}, \eta_{1}\right), V^{*}\left(\xi_{1}, \eta_{1}\right), W^{*}\left(\xi_{1}, \eta_{1}\right)\right),
$$

resulting in a regular curve $\sigma$. Consider the portion $\sigma^{*}$ of $\sigma$ between the two points, since $\sigma^{*}$ is regular and bounded, it must have a $u$-extreme point $p=\left(U^{*}\left(\xi^{*}, \eta^{*}\right), V^{*}\left(\xi^{*}, \eta^{*}\right), W^{*}\left(\xi^{*}, \eta^{*}\right)\right)$ where the normal vector $n$ to the curve is parallel to the $u$-axis, as shown in Fig. 4-1a. Since the projection of the normal $N$ to the surface $X^{*}(\xi, \eta)$ at point $p$ in plane $\Pi$ can be easily seen to identify with $n$, we have $N \cdot w=0$. This translates to

$$
U_{\xi}^{*}\left(\xi^{*}, \eta^{*}\right) V_{\eta}^{*}\left(\xi^{*}, \eta^{*}\right)=U^{*}{ }_{\eta}\left(\xi^{*}, \eta^{*}\right) V_{\xi}^{*}\left(\xi^{*}, \eta^{*}\right)
$$

i.e., $U_{\xi}\left(\xi^{*}, \eta^{*}\right) V_{\eta}\left(\xi^{*}, \eta^{*}\right)=U_{\eta}\left(\xi^{*}, \eta^{*}\right) V_{\xi}\left(\xi^{*}, \eta^{*}\right)$. This means $\left(\xi^{*}, \eta^{*}\right)$ is a singular point of $G(\xi, \eta)$.

Conversely, let $\left(\xi^{*}, \eta^{*}\right)$ be a singular point of $G(\xi, \eta)$; hence, $U_{\xi}\left(\xi^{*}, \eta^{*}\right) V_{\eta}\left(\xi^{*}, \eta^{*}\right)=U_{\eta}\left(\xi^{*}, \eta^{*}\right) V_{\xi}\left(\xi^{*}, \eta^{*}\right)$. Consequently, the normal $N$ to the surface $G^{*}(\xi, \eta)$ at $\left(\xi^{*}, \eta^{*}\right)$ is perpendicular to the $w$-axis. Without loss of generality, we can assume $N$ is parallel to the $u$-axis. Intersecting $G^{*}(\xi, \eta)$ with the plane $v=V^{*}\left(\xi^{*}, \eta^{*}\right)$, we obtain a regular curve $\sigma$. As $p^{*}=\left(U^{*}\left(\xi^{*}, \eta^{*}\right), V^{*}\left(\xi^{*}, \eta^{*}\right), W^{*}\left(\xi^{*}, \eta^{*}\right)\right)$ is a local $u$-extreme point on this curve, one can find a real number $\delta>0$ such that the vertical line $u=U^{*}\left(\xi^{*}, \eta^{*}\right)-\delta$ intersects $\sigma$ at least twice (assuming $p^{*}$ is a $u$-maximum point). Let

$$
p_{0}=\left(U^{*}\left(\xi_{0}, \eta_{0}\right), V^{*}\left(\xi_{0}, \eta_{0}\right), W^{*}\left(\xi_{0}, \eta_{0}\right)\right) \text { and } p_{1}=\left(U^{*}\left(\xi_{1}, \eta_{1}\right), V^{*}\left(\xi_{1}, \eta_{1}\right), W^{*}\left(\xi_{1}, \eta_{1}\right)\right)
$$

be two such intersection points for some $\left(\xi_{0}, \eta_{0}\right) \neq\left(\xi_{1}, \eta_{1}\right)$, as shown in Fig. 4-1b. Obviously, we have $\left(U^{*}\left(\xi_{0}, \eta_{0}\right), V^{*}\left(\xi_{0}, \eta_{0}\right)\right)=\left(U^{*}\left(\xi_{1}, \eta_{1}\right), V^{*}\left(\xi_{1}, \eta_{1}\right)\right)$. Since $(U(\xi, \eta), V(\xi, \eta))=U^{*}(\xi, \eta), V^{*}(\xi, \eta)$, we conclude that $G(\xi, \eta)$ maps two distinct points in the $\xi-\eta$ domain to a same point in the region $\mathrm{P}_{G}$. This completes the proof.
Q.E.D.


Fig. 4-1 Proof of Lemma 4-1.

Based on the above lemma, the following useful proposition is in order.
Proposition 4-1 If the normal at every point to the Gregory patch $G(\xi, \eta)$ has the same sign in $w$, then $G(\xi, \eta)$ has no self-overlapping.

Without loss of generality, we can assume that the sign in $w$ of the normal vectors of a non-self-overlapping mapping $G(\xi, \eta)$ is always positive.

### 4.2 Objective function of the optimization

To facilitate the discussion of the objective function, we use the following terms for self-overlapping.

Definition 4-1 A point $\left(\xi_{d}, \eta_{d}\right)$ is called a shadow point of $G(\xi, \eta)$ if the normal at $G\left(\xi_{d}, \eta_{d}\right)$ is along the negative $w$-axis.

Definition 4-2 The set $R_{d}$ of all the shadow points of $G(\xi, \eta)$ is defined as the shadow region of $G(\xi, \eta)$.

From the above analysis, we find that the non-self-overlapping term in the objective function should be a function indicating the area of the shadow region of $G(\xi, \eta)$. During the algebraic grid generation, the four nodes in every facet are sorted in the counter-clockwise order as $\left(u_{j-1}, v_{j-1}\right),\left(u_{j}, v_{j}\right),\left(u_{j+1}, v_{j+1}\right)$, and $\left(u_{j+2}, v_{j+2}\right)$. Thus, the area of facet $i$ in $M^{G}$ can be computed by

$$
\begin{align*}
& A_{i}=\frac{1}{2}\left[\left(u_{j}-u_{j+1}\right)\left(v_{j}+v_{j+1}\right)+\left(u_{j+1}-u_{j+2}\right)\left(v_{j+1}+v_{j+2}\right)\right.  \tag{8}\\
&\left.\quad+\left(u_{j+2}-u_{j-1}\right)\left(v_{j+2}+v_{j-1}\right)+\left(u_{j-1}-u_{j}\right)\left(v_{j-1}+v_{j}\right)\right]
\end{align*} .
$$

When $A_{i}$ is negative, all the points in facet $i$ are shadow points. Rather than simply adding all the negative $A_{i} \mathrm{~s}$ and taking the sum as the minimization objective function, we adopt an exponential function defined as

$$
\begin{equation*}
J_{1}=\sum_{i=0}^{F-1} e^{-a A_{i}} \tag{9}
\end{equation*}
$$

where $A_{i}$ is the area of facet $i$ in the mesh, $F$ is the total number of facets in the mesh, and $a=\frac{1}{\max \left\{\left|A_{i}\right|\right\}}$. The figure of the function $f(t)=e^{-a t}$ is shown below with $a=2$. The selection of this function over simple summation is motivated by considerations: (1) this function usually converges better than the simple linear summation in the iterative minimization process, and (2) it also tends to compute a global minimization. Why not simply applying a spring model (i.e., $f(t)=\frac{1}{2} a t^{2}$ ) as the format of our objective function? It is because that we want the first term $J_{1}$ effecting on the grids only when having some negative $A_{i}$ s. Thus, the ideal format should have $f(t) \approx 0$ when $t>0$. Also, the function $f(t)$ is expected to be $C^{1}$ continuous since our objective function will be minimized by a gradient method. The exponential function satisfies these factors but a spring model does not.


Fig. 4-2 Shape of function $f(t)=e^{-a t}$ with $a=2$

To achieve a smooth grid with good facet shape, Laplacian smoothing is usually applied on structured grids, which arises from solving a pair of partial differential equations [19]: $u_{\xi \xi}+u_{\eta \eta}=0$ and $v_{\xi \xi}+v_{\eta \eta}=0$. In practice, node positions are the average of points of its neighboring nodes in Laplacian smoothing. Similarly, we add the following smoothing term into our objective function

$$
\begin{equation*}
J_{2}=\sum_{j=0}^{E-1}\left\|\vec{v}_{j}-\frac{1}{L} \sum_{m \in j^{*}} \vec{v}_{m}\right\|^{2} \tag{10}
\end{equation*}
$$

where $j^{*}$ represents an index complex of the neighbor nodes of the inner node $\vec{v}_{j}=\left(u_{j}, v_{j}\right)$ in $M^{G}, L$ is the element number of $j^{*}$, and $E$ is the number of inner nodes in the mesh. The formula of $J_{2}$ can be rewritten as

$$
\begin{equation*}
J_{2}=\sum_{j=0}^{E-1}\left[\left(u_{j}-\frac{1}{L} \sum_{m \in j^{*}} u_{m}\right)^{2}+\left(v_{j}-\frac{1}{L} \sum_{m \in j^{*}} v_{m}\right)^{2}\right] . \tag{11}
\end{equation*}
$$

By summing together the non-self-overlapping term and the smoothing term, the final objective function is given as

$$
\begin{equation*}
J=\sum_{i=0}^{F-1} e^{-a A_{i}}+\frac{1}{\bar{J}_{2}} \sum_{j=0}^{E-1}\left[\left(u_{j}-\frac{1}{L} \sum_{m \in j^{*}} u_{m}\right)^{2}+\left(v_{j}-\frac{1}{L} \sum_{m \in j^{*}} v_{m}\right)^{2}\right], \tag{12}
\end{equation*}
$$

where $\bar{J}_{2}$ is the initial value of $J_{2}$ before the optimization. This factor is utilized to balance the relative importance of the non-self-overlapping term and the smoothing term so that $J_{1}$ leads the nodes' movement when self-overlapping occurs, and $J_{2}$ directs the nodes to archive better grid shape when $J_{1} \rightarrow 0$.

### 4.3 Numerical implementation

During the process of optimization, we move the inner nodes of $M^{G}$ in the $u-v$ plane to achieve the functional optimum (minimum). Therefore, the $u$ and $v$ components of all the inner nodes of $M^{G}$ form the solution vector $\chi$. The Conjugate Gradient method [20] is applied to obtain the functional optimum, where the explicit form of gradient at every inner node to the objective function is desired. For an inner node $\left(u_{j}, v_{j}\right)$, from equation (8), we have

$$
\frac{\partial A_{i}}{\partial u_{j}}=\frac{1}{2}\left(v_{j+1}-v_{j-1}\right) \text { and } \frac{\partial A_{i}}{\partial v_{j}}=\frac{1}{2}\left(u_{j-1}-u_{j+1}\right),
$$

so the gradient direction of $\left(u_{j}, v_{j}\right)$ with respect to the non-self-overlapping term $J_{1}$ in the objective function is

$$
\left[\begin{array}{l}
\frac{\partial J_{1}}{\partial u_{j}}  \tag{13}\\
\frac{\partial J_{1}}{\partial v_{j}}
\end{array}\right]=\left[\begin{array}{l}
\sum_{k=0}^{N-1}\left(-a e^{-a A_{k}}\right) \frac{\partial A_{k}}{\partial u_{j}} \\
\sum_{k=0}^{N-1}\left(-a e^{-a A_{k}}\right) \frac{\partial A_{k}}{\partial v_{j}}
\end{array}\right]=\left[\begin{array}{l}
\frac{1}{2} \sum_{k=0}^{N-1}-a e^{-a A_{k}}\left(v_{j+1}-v_{j-1}\right) \\
\frac{1}{2} \sum_{k=0}^{N-1}-a e^{-a A_{k}}\left(u_{j-1}-u_{j+1}\right)
\end{array}\right],
$$

where $A_{k}$ is the area of the $k$ th incident facet of $\left(u_{j}, v_{j}\right)$, and $N$ is the number of incident facets around $\left(u_{j}, v_{j}\right)$. By equation (11), the gradient of $\left(u_{j}, v_{j}\right)$ with respect to the smoothing term $J_{2}$ in the objective function is determined as

$$
\left[\begin{array}{l}
\frac{\partial J_{2}}{\partial u_{j}}  \tag{14}\\
\frac{\partial J_{2}}{\partial v_{j}}
\end{array}\right]=\left[\begin{array}{l}
2\left(u_{j}-\frac{1}{L} \sum_{m \in j^{*}} u_{m}\right)-\sum_{m \in j^{*}} \frac{2}{L_{m}}\left(u_{m}-\frac{1}{L_{m}} \sum_{n \in m^{*}} u_{n}\right) \\
2\left(v_{j}-\frac{1}{L} \sum_{m \in j^{*}} v_{m}\right)-\sum_{m \in j^{*}} \frac{2}{L_{m}}\left(v_{m}-\frac{1}{L_{m}} \sum_{n \in m^{*}} v_{n}\right)
\end{array}\right],
$$

where $m^{*}$ represents an index complex of the neighboring nodes of $\left(u_{j}, v_{j}\right)$, and $L_{m}$ is the element number of $m$ *. Therefore, in summary, we have

$$
\left[\begin{array}{l}
\frac{\partial J}{\partial u_{j}}  \tag{15}\\
\frac{\partial J}{\partial v_{j}}
\end{array}\right]=\left[\begin{array}{l}
\frac{1}{2} \sum_{k=0}^{N-1}-a e^{-a A_{k}}\left(v_{j+1}-v_{j-1}\right)+\frac{2}{\bar{J}_{2}}\left[\left(u_{j}-\frac{1}{L} \sum_{m \in j^{*}} u_{m}\right)-\sum_{m \in j^{*}} \frac{1}{L_{m}}\left(u_{m}-\frac{1}{L_{m}} \sum_{n \in m^{*}} u_{n}\right)\right] \\
\frac{1}{2} \sum_{k=0}^{N-1}-a e^{-a A_{k}}\left(u_{j-1}-u_{j+1}\right)+\frac{2}{\bar{J}_{2}}\left[\left(v_{j}-\frac{1}{L} \sum_{m \in j^{*}} v_{m}\right)-\sum_{m \in j^{*}} \frac{1}{L_{m}}\left(v_{m}-\frac{1}{L_{m}} \sum_{n \in m^{*}} v_{n}\right)\right]
\end{array}\right] .
$$

Using the above formula, by a Conjugate Gradient method, we can iteratively determine the value of components in $\chi$ that makes the objective function as defined in equation (12) minimum. Thus, the $u$ and $v$
parameters of every inner node in a non-self-overlapping grid $M^{G^{*}}$ are determined. By the parametric equation of $S_{T}$, the final position of all grid nodes can be computed by their $u, v$ parameters.


Fig. 4-3 Example II - grid before vs. after optimization

Fig. 4-3 shows an example of the grid (in $u-v$ plane) generated before and after the functional optimization when choosing a $10 \times 10$ grid in each sub-domain. As demonstrated in the figure, the selfoverlapping is eliminated in the optimized grid.

### 4.4 Progressive optimization

As our objective function in the functional optimization is concave with respect to its solution vector $\chi$ (there can be many local minima), the success of the numerical optimization algorithm depends critically on the initial position of $\chi$. We take the algebraic grid generated by a Gregory patch in section 3 as an initial position of the solution vector; it however may not be a very satisfying one for some strongly concaved boundaries (e.g., the one shown in Fig. 4-4a). "Guessing" a good initial vector is hard. The basic idea we take to overcome this difficulty is to progressively achieve the optimum by gradually deforming the grid $M^{0}$ from a regular $n$-gon $\Omega_{R}$ into the grid $M^{f}$ in the region $\mathrm{P}_{S_{T}}$ bounded by the given curves in the $u-v$ plane, which we discuss in detail in this section.

The grid $M^{0}$ generated in the initial shape, a regular $n$-gon, by a Gregory patch $G^{0}(\xi, \eta)$ is non-selfoverlapping. For any grid node $\vec{v}_{j}$, its coordinates determined by the Gregory patch $G^{T}(\xi, \eta)$ on $\mathrm{P}_{S_{T}}$ in the $u-v$ plane are $\left(u_{j}^{T}, v_{j}^{T}\right)$, and the coordinates determined by $G^{0}(\xi, \eta)$ are $\left(u_{j}^{0}, v_{j}^{0}\right)$. At the beginning of the deformation, the moving direction of $\vec{v}_{j}$ is $\left(d u_{j}, d v_{j}\right)=\left(u_{j}^{T}-u_{j}^{0}, v_{j}^{T}-v_{j}^{0}\right)$. When deforming $M^{0}$ into $M^{f}$ by changing a deformation factor $\lambda_{t}\left(\lambda_{t} \in[0,1]\right)$, the position of every grid node $\vec{v}_{j}$ is determined by

$$
\begin{align*}
u_{j}^{t} & =u_{j}^{0}+\lambda_{t} d u_{j}  \tag{16}\\
v_{j}^{t} & =v_{j}^{0}+\lambda_{t} d v_{j}
\end{align*}
$$

During the deformation, we check if self-overlapping occurs; once it is detected, the numerical optimization method presented in section 4-3 is applied to current grid $M^{\lambda_{t}}$ to construct a new grid $M^{\lambda_{t}{ }^{*}}$ without or with less self-overlapping. The position of grid node $\vec{v}_{j}$ in $M^{\lambda_{t}{ }^{*}}$ changes from $\left(u_{j}^{t}, v_{j}^{t}\right)$ to $\left(u_{j}^{t^{*}}, v_{j}^{t^{*}}\right)$, so we use the following equation to alternate the moving direction of $\vec{v}_{j}$.

$$
\begin{align*}
d u_{j} & =\left(u_{j}^{t^{*}}-u_{j}^{0}\right) / \lambda_{t}  \tag{17}\\
d v_{j} & =\left(v_{j}^{t_{j}^{*}}-v_{j}^{0}\right) / \lambda_{t}
\end{align*} .
$$

The deformation will then continue along the new directions guided by equation (16). The deformation and functional optimization are applied alternatively until the final non-self-overlapping grid $M^{f}$ is obtained. Since the positions of boundary grid nodes are not adjusted during the optimization, their moving directions do not change during the deformation - so the result grid $M^{f}$ of the progressive optimization is still boundary conforming to $\mathrm{P}_{S_{T}}$ exactly as the algebraic grid $M^{T}$ generated by the Gregory patch $G^{T}(\xi, \eta)$ on $\mathrm{P}_{S_{T}}$.

During the deformation, the deformation factor $\lambda_{t}$ increases from zero to one adaptively to the value increase of the objective function. The overall procedure of progressive optimization is outlined in pseudo-code in Algorithm ProgressiveOptimization() below. As a demonstration of the power of this progressive optimization, to Example III (shown in Fig. 4-4), we first tried the pure numerical optimization which though failed to achieve a non-self-overlapping grid even after iterating 10000 times; using Algorithm ProgressiveOptimization(), we easily obtain a final result without self-overlapping by applying the numerical optimization only 8 times with summed 479 iterations. The progressive results are shown in Fig. 4-5. The progressive algorithm does not success in any case, so in step 7 of Algorithm ProgressiveOptimization() we detect whether the moving step is less than a threshold. If so, report failure and stop running the algorithm. However, in all of our testing examples, the algorithm works well.

Algorithm ProgressiveOptimization ( $\mathrm{P}_{S_{T}}$ )
Input: A region $\mathrm{P}_{S_{T}}$ in the $u-v$ parametric space.
Output: The non-self-overlapping grid $M^{f}$ on $\mathrm{P}_{S_{T}}$.

1. Compute the algebraic grid $M^{T}$ on $\mathrm{P}_{S_{T}}, M^{T}$ has $n_{T}$ facets;
2. Compute position of every grid node on $M^{0}$;
3. Compute the initial moving direction of every grid node by $G^{0}(\xi, \eta)$ and $G^{T}(\xi, \eta)$;
4. $\lambda_{t} \leftarrow 0$;
5. do $\{$
6. $\Delta \lambda \leftarrow 2\left(1-\lambda_{t}\right)$;
7. When $\Delta \lambda<10^{-16}$, report failure and terminate the algorithm;
8. do \{
9. $\Delta \lambda \leftarrow \Delta \lambda / 2$, and $\lambda_{t} \leftarrow \lambda_{t}+\Delta \lambda$;
10. Change the position of every grid node by equation (16) - obtain the current grid $M^{\lambda_{t}}$;
11. Compute the number $n_{-}$of facets with negative area on the current grid $M^{\lambda_{t}}$;
12. $\}$ while $\left(\left(n_{-} / n_{T}\right)>\tau\right)$;
13. Compute the numerical optimum $M^{\lambda_{t}{ }^{*}}$ of the current grid $M^{\lambda_{t}}$;
14. Use $M^{\lambda_{t} *}$ to update the moving direction of every grid node by equation (17);
15. $\}$ while $\left(\lambda_{t}<1\right)$;
16. $M^{f} \leftarrow M^{\lambda_{t}{ }^{*}}$;
17. return $M^{f}$;
(* in our testing, we choose $\tau=20 \%$ )

(a) before optimization

(b) after optimization

Fig. 4-4 Example III - result of progressive optimization


Fig. 4-5 Example III - progressive results

## 5. Experimental Results

The proposed mesh algorithm has been implemented using Java language and separately tested on a PIII 900 MHz PC with a basic configuration and on a PIV 2.6 GHz PC with a modernist configuration; a number of test cases are tried. Fig. 5-1 gives the non-self-overlapping grid generation result of the trimmed surface in Example I (initially given in Fig. 1-2); Fig. 5-2 shows the result of the patch in Example II with a mesh denser than the one shown in Fig. 4-3; and the final surface grids in Example III are displayed in Fig. 5-3. Example IV is a trimmed surface with six sides; its shape in the parametric surface is very convoluted so that the 6 -side region cannot be easily divided into two 4 -side sub-regions by straight lines since all the pertinent diagonals (shown as dash lines in Fig. 5-4a) intersect the boundaries. As shown in Fig. 5-4b, the original algebraic grids from the Gregory patch mapping incurs severe self-overlapping. Using the presented mesh method, the selfoverlapping is successfully eliminated in the final optimized grid, as given in Fig. 5-4c. The final surface grids are shown in Fig. 5-4d. For a trimmed surface patch with holes, we can first partition its parametric domain into several regions without holes and then apply the proposed mesh algorithm independently to these regions, e.g.

Example V shown in Fig 5-5. All the boundary curves in our examples are presented by $4^{\text {th }}$ order Bezier curves. The control points of the Bezier boundary curves are listed in Table 5-1.

The computer running times of the given examples are tabulated in Table 5-2. Usually, several minutes are needed. If the standard tools in commercial software are conducted to generate the grids on trimmed $n$-sided parametric patches, manual partitioning is required and it is a trial and error process - so it usually takes several hours of a human working. Also, it is hard to control the time cost. In some cases, especially the cases with convex boundaries, the traditional process can generate a good result very fast. However, if you are unluckily dealing with a much concaved parametric region, you may spend a lot of time on it. It is believed that with the increasing processing power available on the computer and with more efficient optimization algorithms, the running time can be shortened significantly. Consider about the differences of computing time on PCs with basic and modernist configurations, the computing time of all examples on PIV 2.6 GHz is almost one third of the time on PIII 900 . On the PIV 2.6 GHz , it has already achieved an acceptable speed in the work of computeraided engineering.


Fig. 5-1 Result of Example I

(a) result grids in $u-v$ space

(b) final grids on the trimmed surface

Fig. 5-2 Result of Example II


Fig. 5-3 Result of Example III

(a) the parametric area in $u-v$ domain
(c) grids after progressive optimization


(b) algebraic grids in $u-v$ domain

Fig. 5-4 Example IV - a patch with six sides


Fig. 5-5 Example V - complex region with multiple patches

Table 5-2 Time cost of examples

| Example | Result <br> Figures | Optimization Type | Side <br> Number* | Time cost on <br> PIII 900MHz | Time cost on <br> PIV 2.6GHz |
| :---: | :---: | :---: | :---: | :---: | :---: |
| I | $5-1$ | Progressive | 5 | 4 min .30 sec. | 1 min .41 sec. |
| II | $4-3,5-2$ | Pure numerical | 5 | 2 min .11 sec. | 49 sec. |
| III | $4-4,4-5,5-3$ | Progressive | 5 | 12 min .7 sec. | 4 min .13 sec. |
| IV | $5-4$ | Progressive | 6 | 15 min .13 sec. | 5 min .19 sec. |
| V | $5-5$ | Progressive | $2 \times 5$ | 4 min .9 sec. | 1 min .33 sec. |

[^1]Table 5-1 Control points of boundary curves

| Example | Curve No. | Control Points |
| :---: | :---: | :---: |
| I | 1 | (0.86,0.90), (1.0,0.65), (0.67,0.40), (0.41,0.0) |
|  | 2 | (0.41,0.0), (0.30,0.08), (0.26,0.01), (0.0,0.21) |
|  | 3 | $(0.0,0.21),(0.53,0.31),(0.56,0.42),(0.53,0.52)$ |
|  | 4 | (0.53, 0.52$),(0.73,0.61),(0.60,0.79),(0.27,0.82)$ |
|  | 5 | (0.27,0.82), (0.34,0.96), (0.63,1.0), (0.86,0.90) |
| II | 1 | (0.93, 0.46$),(0.88,0.33),(0.89,0.14),(0.65,0.01)$ |
|  | 2 | (0.65,0.01), (0.38,0.0), (0.42,0.20), (0.10,0.20) |
|  | 3 | (0.10,0.20), (0.72,0.40), (0.46,0.57), (0.0,0.44) |
|  | 4 | (0.0,0.44), (0.26,0.87), (0.46,0.94), (1.0,0.99) |
|  | 5 | (1.0,0.99), (0.32,0.54), (0.66,0.29), (0.93,0.46) |
| III | 1 | $(1.0,0.36),(0.50,0.37),(0.26,0.18),(0.92,0.10)$ |
|  | 2 | (0.92,0.10), (0.46,0.0), (0.21,0.10), (0.0,0.30) |
|  | 3 | (0.0,0.30), (0.94,0.24), (0.67,1.0), (0.15,0.34) |
|  | 4 | (0.15,0.34), (0.12,0.44), (0.15,0.69), (0.31,0.79) |
|  | 5 | (0.31,0.79), (0.70,0.70), (0.82,0.59), (1.0,0.36) |
| IV | 1 | (0.62,0.40), (0.74,0.38), (1.0,0.31), (0.90,0.10) |
|  | 2 | (0.90,0.10), (0.73,0.21), (0.58,0.23), (0.27,0.0) |
|  | 3 | (0.27,0.0), (0.37,0.29), (0.28,0.65), (0.0,0.64) |
|  | 4 | (0.0,0.64), (0.12,0.86), (0.37,1.0), (0.46,0.75) |
|  | 5 | (0.46,0.75), (0.62,0.74), (0.77,0.66), (0.84,0.53) |
|  | 6 | (0.84,0.53), (0.71,0.52), (0.59,0.49), (0.62,0.40) |

## 6. Conclusion

In this paper, we present a method for automatically constructing a structured grid system in an $n$-sided planar region bounded by parametric boundary curves of any form with only $C^{1}$ continuity, using a planar Gregory patch. This averts the need of manually partitioning the $n$-sided region into 4 -sided sub-regions, which is a popular solution by most commercial meshing software. However, the algebraic grid thus generated may have self-overlapping which makes the mesh useless in most of engineering and scientific applications. A functional optimization method is then presented to eliminate such self-overlapping in the grid. Unlike PDE
methods, the derivatives of a Gregory patch, which are very difficult to compute, are not required in our method. Thus, the speed of the optimization process of our approach is relatively fast. To resolve the difficulty of guessing good initial positions of every grid node for the conjugate gradient method, a progressive optimization algorithm is introduced, which has been shown to be very effective in a variety of practical examples. In summary, our approach provides a promising meshing tool in engineering.

Some new ideas about introducing additional terms to the Gregory definition have been proposed recently [18], which will allow extra controlling to the patch. One possible extension of our current work is to see if an algebraic mapping which guarantees non-self-overlapping can be determined by adjusting those additional terms of a Gregory patch. The process may also be a functional optimization approach like what is described in this paper. Or we can alternate the 3D vector functions along the boundaries so that the positions of inner nodes are adaptively modified to achieve a non-self-overlapping grid.

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[^1]:    *The grid size in each block is $20 \times 20$.

