# Universidade Federal de Juiz de Fora <br> Instituto de ciências Exatas <br> Bacharelado em Ciência da Computação 

# Iterative method for edge equalization of triangular meshes 

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#### Abstract

MONOGRAFIA SUBMETIDA AO CORPO DOCENTE DO INSTITUTO DE CIÊNCIAS EXATAS DA UNIVERSIDADE FEDERAL DE JUIZ DE FORA, COMO PARTE INTEgrante dos requisitos necessários para a obtenção do grau de BaCHAREL EM CIÊNCIA DA COMPUTAÇÃO.


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To my friends and my sister.
To my Parents for their support and sustenance.

## Resumo

Este trabalho apresenta um método para remalhamento de superfícies triangulares. O método obtem uma nova superfície de maneira que qualquer aresta está dentro de um intervalo predefinido $\left[e_{\min }, e_{\max }\right]$. A entrada do processo é uma malha de 2-variedade com geometria e topologia arbitrarias. O algoritmo proposto é iterativo e consegue ajustar automaticamente a quantidade de vértices e triângulos necessários através das operações estelares. Um filtro passa-baixa também é aplicado para retirar as altas frequências. O algoritmo gera uma malha triangular de 2-variedade, com os vértices distribuídos de maneira quase uniforme sobre a superfície de entrada. No fim do processo, praticamente todas as arestas estão dentro do intervalo. A malha dual desta malha triangular é uma malha trivalente. Este tipo de malha tem muitas aplicações em simulações de nano estruturas de carbono.

Palavras-chave: equalização do comprimento de aresta, operações estelares, remalhamento.


#### Abstract

This work presents a method for remeshing triangular surfaces. This method obtains a new mesh in such a way that any edge is within a predefined interval $\left[e_{\min }, e_{\max }\right]$. The input of the process is a 2-manifold mesh with arbitrary geometry and topology. The proposed algorithm is iterative and is able to automatically adjust the number of vertices and polygons through stellar operations. A low pass filter is also applied to remove higher frequencies. This algorithm generates a triangular 2-manifold mesh, with the vertices spreaded almost uniform over the input surface. At the end of the process, almost all edges are within the interval. The dual mesh of this triangular mesh is a trivalent mesh also very uniform. This kind of mesh has many applications in simulations of carbon nano structures.


Keywords: edge length equalization, stellar operations, remesh.

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## 1 Introduction

This work aims the production of a method capable of, from a given mesh, generate a new mesh with edge lengths that are within a minimum and maximum value.

To understand this work it is necessary that some basic concepts are clear. These concepts are: polygon meshes to represent surfaces, triangular meshes, stellar operations, the laplacian operator and remeshing. To represent a surface in the computer memory a discretization is necessary. To achieve this there are several manners. One of the most used is to represent objects through polygon meshes. Polygon meshes are surfaces represented by discrete polygons, each polygon representing the surface locally and the entire surface is represented by all polygons together.

The polygon meshes can be made with different types of polygons, for example triangles, hexagons, quads, pentagons. In this work we are going to focus only in triangular meshes due to their great utility and practical use. Other very important mesh in this work are the trivalent meshes, in which each vertex has exactly three edges. They tend to form hexagons in low curvature surfaces and have recently attracted attention from the computer geometry community (M. Nieser et al, 2010). They are important for several applications, such as the physics simulation of carbon nano structures (Quinelato et al, 2010).

Stellar operations are a classic set of operations in polygon meshes known for keeping the Euler characteristic unchanged as in equation 1.1

$$
\begin{equation*}
v_{b}+e_{b}+f_{b}=v_{a}+e_{a}+f_{a}, \tag{1.1}
\end{equation*}
$$

where $v_{b}$ is the original number of vertices, $e_{b}$ is the original number of edges, $f_{b}$ is the original number of faces, $v_{a}$ is the new number of vertices, $e_{a}$ is the new number of edges, and $f_{a}$ is the new number of faces. They are formed basically by the operations of edge split, vertex split, edge collapse, vertex collapse and edge flip (Oliveira et al, 2012).

The Laplacian operator is a common operator often used as a filter, as it tends to
remove the higher frequencies. Basically it is an operator that try to reposition the vertex in the centroid of its neighbors. In this work, we use the laplacian to act not only in the first neighbors but in the $k$-neighborhood with a differential weight to each neighborhood (Oliveira et al, 2012).

According to Mario et al (2010a), remeshing is an operation that generates a new mesh with greater quality than its original version. The term quality is dependent on the objective. For example, to minimize memory, the objective can be to reduce the number of polygons and vertices without losing accuracy. For this work, a quality mesh is the one that contains all edge lengths in the target interval. This is extremely useful in atomic simulations where the physically plausible distances between the atoms are known and limited.

### 1.1 Problem definition

Given an arbitrary triangular mesh $\mathcal{M}$, composed of edges $E_{i} \in \mathcal{M}$, generate a new trivalent mesh with the same geometry and topology, but having all edge lengths in the interval $\left[e_{\min }, e_{\max }\right]$. In other words, $e_{\min } \leq\left|E_{i}\right| \leq e_{\max } \forall E_{i} \subset \mathcal{M}$.

### 1.2 Motivation

In this work, we propose a new iterative process to transform a triangular mesh in a trivalent mesh whose edge lengths are within a predefined interval. The method was conceived to be used with most 2-manifold models found publicly.

In a previous work (Oliveira et al, 2012),(Oliveira et al, 2013), an iterative method was developed to force the median and standard-deviation of the mesh to be below expected thresholds. This works has the drawback of allowing some edges to be much greater than the overall average. The imposition of a minimum and maximum value for edge lengths is much more restrictive since the majority of the edges cannot compensate the extra size of a few outliers. This problem statement constrains even further the space of possible solutions for the mesh, and is a motivator for this work.

### 1.3 Objectives

The main objective of this work is to develop a method to transform a triangular mesh in a trivalent mesh with all edges within an interval. This method must preserve the geometry and topology of the surface.

As a secondary objective, the method also try to obtain a hexagonal mesh, in order to be practical for nano structure simulations. Another objective is to have low memory cost since some input meshes might have hundreds of thousands vertices. Methods based on linear system solving require large amounts of installed memory to process big meshes. We propose a iterative method capable of processing the large meshes found publicly.

## 2 Basics Concepts

### 2.1 Manifolds

A manifold represents a $n$ dimension surface immersed in $k, k \geq n$ dimension space. In this kind of surfaces the neighborhood of each point can be reduced to an $n$ dimension euclidean space (Guillemin and Pollack, 1974).

A surface is orientable if it has two sides. As strange as it can be, there are some surfaces that do not have two sides. The Möbius strip is a classic example and is depicted in Figure 2.1.


Figura 2.1: This is an example of non orientable surface (Starostin et al, 2010).

This work will be focused in surfaces that are orientable.

### 2.2 Euler characteristic

The Euler characteristic is a topological invariant of surfaces. In other words, any two surfaces that are homeomorphic will have the same Euler characteristic, it does not matter their size, shape or any other characteristic of the surface.

The geometric methods of this work must preserve the Euler characteristic in every operation to consequently preserve the topology of the surface.

### 2.3 Triangular meshes

A surface could be represented in many manners, a possible discretization with great use in computer graphics, due this simplicity and versatility are the triangular meshes. This is a especial kind of mesh where the surface is approximated through triangles, represented by their distinct and non-collinear three vertices $[A, B, C]$. Due to the barycentric parameterization, each point $P$ in a triangle $[A, B, C]$ is represented as an unique linear combination:

$$
\begin{array}{r}
P=\alpha A+\beta B+\gamma C, \\
\alpha+\beta+\gamma=1 . \tag{2.2}
\end{array}
$$

Triangulation is an efficient way for approaching every point over the surface with a very simple data structure. Besides the simplicity, this approach has a great power of representation. Any three distinct and non-collinear points always generate a plan, which is simpler to deal geometrically.

The vertex valency is the number of incident edges. An equilateral triangle has internal angles equals to sixty degrees. In a scene of ideal valency for the purposes of this work, each vertex should have a valency of $\frac{360^{\circ}}{60^{\circ}}=6$. In this work, a vertex with valency six is called a regular vertex and all other vertex valencies are called irregular vertex (Oliveira et al, 2012).

In this work, all input meshes will be implicitly orientable and represented by a triangular mesh.

### 2.4 Stellar Operations

Stellar Operations are operations that preserve the Euler characteristic of the mesh (Mario et al, 2010a), they are based on division, remotion or changing edges from the model however maintaining the sum vertices - edges + faces of the surface as showed in Equation 1.1. Consider the vertices $V_{1}, V_{2}, V_{3}$ and $V_{4}$ forming two triangles $f_{1}=\left(V_{1}, V_{2}, V_{3}\right)$ and $f_{2}=\left(V_{1}, V_{3}, V_{4}\right)$. The vertices adjacent to the edge $A_{i}=\overline{V_{1} V_{3}}$ are $V_{1}$ and $V_{3}$, the opposite
vertices to $A_{i}$ are $V_{2}$ and $V_{4}$. The set of vertices from the triangles adjacent to $A_{i}$ is $\left\{V_{1}, V_{2}, V_{3}, V_{4}\right\}$

### 2.4.1 Edge flip

This operation flips an edge as depicted in Figure 2.2.


Figura 2.2: Edge Flip (Oliveira et al, 2012).

In this operation the edge $A_{i}$ is changed by a new edge linking the opposite vertices of $A_{i}$, the Euler characteristic is preserved.

### 2.4.2 Edge split

There are four vertices involved in a edge split. After the split, a new vertex is created over the split edge, two new faces will be added in the mesh. The valency of the opposite vertices will be increased by one as depicted in Figure 2.3.


Figura 2.3: Edge Split (Oliveira et al, 2012).

### 2.4.3 Edge collapse

This operation removes an edge from the mesh as depicted in Figure 2.4. There are four


Figura 2.4: Edge collapse (Oliveira et al, 2012).
vertices involved in a collapse. An edge $A_{i}$ is selected and removed, the faces containing the edge $A_{i}$ are removed too. Given the vertices that form the edge $\overline{V_{i} V_{j}}$, only one is maintained and its valency $v a_{i}$ becomes $v a_{i}+v a_{j}-4$.

### 2.5 Remesh

According to Mario et al (2010a), a remesh is defined as: Given a surface represented by a mesh $\mathcal{M}$, create a new mesh $\mathcal{M}^{\prime}$, such that the new mesh satisfy a certain quality standard and $\mathcal{M}^{\prime}$ is enough next of $\mathcal{M}$. So remesh is a very important technique when you a have a certain standard to be achieved by a surface. In this case, the goal is to have the edges length within a predefined interval. The definition of "enough next" is vague. It could be topological, geometric or defined by any constraint. In general, the topology must be preserved while other characteristics might be approximately the same as the original (Surazhsky et al, 2003).

### 2.6 Lowpass filtering

A lowpass filter removes high frequencies from the mesh, making it more smooth. There are several lowpass filters and, in this work, we use a Laplacian filter. More precisely, we use a modified Laplacian filter which uses more than one ring around the vertices ( $k$-rings) and is constrained to be in a previously defined tangent plan.

The classic laplacian filter is defined as:

$$
\nabla^{2} f=\frac{\partial^{2} f}{\partial^{2} x_{1}}+\ldots+\frac{\partial^{2} f}{\partial^{2} x_{n}}
$$

It is a measure of the dispersion in $\mathbb{R}^{n}$ of a function $f$. Taubin et al (1995) propose a discrete approach to the Laplacian operator. The approach is:

$$
\begin{equation*}
L\left(V_{i}\right)=\sum_{V_{j}} w_{i j}\left(V_{i}-V_{j}\right), \tag{2.3}
\end{equation*}
$$

with $V_{j}$ in the neighborhood of $V_{i}$. In the literature, many weights were proposed for $w_{i j}$. There are schemes based on cotangent (Alliez et al, 2002) and neighborhood (Oliveira et al, 2012). The discrete Laplacian is largely used due to its simplicity. Basically, each vertex is moved to the median of its neighbors. This procedure tends to equalize edges lengths, minimizing the standard deviation. The Laplacian must be zero to achieve these properties and the system to solved is given by:

$$
\sum_{v_{j}} w_{i j}\left(v_{i}-v_{j}\right)=0 .
$$

In this work, we use an iterative approach that approximately gives the same results of the Laplacian filter.

## 3 Proposed method

The input for the algorithm that uniforms the edges is a tuple ( $\left.\mathcal{M}, e_{\text {min }}, e_{\text {max }}, k, n, p r, f\right)$, where $\mathcal{M}$ is the triangular mesh, $e_{\text {min }}$ is the smallest edge length allowed, $e_{m} a x$ is the biggest edge length allowed, $k$ the number of rings used at the Laplacian optimization step, $n$ the number of iterations, $p r$ is the number of iterations before the original mesh is replaced by the current mesh in order to relax next projections, $f$ is the maximum degradation for the triangles smallest angle involved in the flip. The loss of the original geometry is smaller as $f$ is closer to zero.

```
Algoritmo 1: UniformRemeshing( \(\left.\mathcal{M}, e_{\text {min }}, e_{\text {max }}, k, n, p r, f\right)\)
    \(\mathcal{M}^{\prime}=\operatorname{Copy}(\mathcal{M})\)
    \(m=\frac{e_{\min }+e_{\text {max }}}{2}\)
    for \(i=1\) to \(n\) do
        if \(p r>0\) and \((\) imodpr \()=0\) then
            \(\mathcal{M}=\operatorname{Copy}\left(\mathcal{M}^{\prime}\right)\)
        end if
        \(m_{i}=\operatorname{MIN}\left(2 \cdot\right.\) CalculateEdgesAverage ( \(\left.\mathcal{M}^{\prime}\right), m\) )
        \(L_{p}=\) CreatePriorityList ( \(\mathcal{M}^{\prime}, m_{i}, e_{\text {min }}, e_{\text {max }}\) )
        StellarOperations ( \(\mathcal{M}^{\prime}, L_{p}\) )
        CorrectValency ( \(\mathcal{M}^{\prime}\) )
        LowPassFiltering ( \(\left.\mathcal{M}^{\prime}, k\right)\)
        Projection( \(\mathcal{M}, \mathcal{M}^{\prime}\) )
    end for
    return \(\mathcal{M}^{\prime}\)
```

The method works as shown in Algorithm 1. Each step is separately explained ahead.

### 3.1 Stellar Transformations with Priority List

The mesh must be transformed according to the target median $m=\frac{e_{\text {min }}+e_{\text {max }}}{2}$. Less faces and vertices are needed to represent the model as $m$ is higher. For the opposite, a lower $m$ results in more faces and vertices.

The current edge's length is the criterion to decide if vertices are added or re-
moved. An arbitrary input mesh, however, can have regions which should be refined and others that should be simplified.

This step is based on the modification of long or short edges $A_{j}$ which are classified as:

$$
\begin{aligned}
\text { long, } & \text { if }\left|A_{j}\right|>m_{i}+\sigma, \\
\text { short, } & \text { if }\left|A_{j}\right|<m_{i}-\sigma,
\end{aligned}
$$

where $\sigma=\frac{e_{\text {max }}-e_{\text {min }}}{2}$ and $m_{i}$ is an intermediate target value for the $i$-th iteration, defined for the transition between the iteration $i$ and $i+1$. This intermediate target value $m_{i}$ avoids sudden changes on the mesh geometry if $m$ is too different than the original mesh edge length average.

Edges with lengths within the closed interval $\left[m_{i}-\sigma, m_{i}+\sigma\right]$ remain unchanged during this step. Long and short edges are then candidates to be collapsed or split, respectively. For every edge $A_{i} \in \mathcal{M}^{\prime}$ considered long, a new vertex is inserted over the edge, dividing it into two smaller edges by the stellar operation of edge split. Edges considered too small in respect to $m_{i}$ are removed from the model using edge collapse.

When $m$ is much greater than the current edges average, a strong mesh simplification is required. In a extreme example with high $m$, all edges will be candidates to be collapsed. Thus, the iteration's target $m_{i}$ is set to be the minimum between the global target average $m$ and two times the current edges average of $\mathcal{M}_{i}^{\prime}$ (Algorithm 1). In that way, the difference between the averages of $\mathcal{M}_{i}^{\prime}$ and $\mathcal{M}_{i+1}^{\prime}$ changes smoothly and is unlikely to generate invalid triangles or change the topology of the mesh, due to the collapse of near long edges. The upper bound of two times the current average per iteration for $m_{i}$ comes from the fact that the edges progressively tend to be equalized, and doubling the edges length of a star does not promote a severe local modification.

The order of application of stellar operations is important. We propose to process the longest and smallest edges first. A priority list $L_{p}$ is created where the edges $A_{j}$, with higher deviation $d_{j}=\left|\left|A_{j}\right|-m_{i}\right|$, are the most important. The list $L_{p}$ is the set of edges $\left\{A_{1}, \ldots, A_{t}\right\} \subset \mathcal{M}$, with $d_{1} \geq d_{2}, \ldots, d_{t-1} \geq d_{t}$, where $A_{j} \in L_{p}$ if $\left|A_{j}\right| \notin\left[m_{i}-\sigma, m_{i}+\sigma\right] \in \mathbb{R}$ (Algorithm 2).

The stellar operations are performed after the set up of the list. The algorithm

```
Algoritmo 2: CreatePriorityList( \(\left.\mathcal{M}^{\prime}, m_{i}, e_{\text {min }}, e_{\text {max }}\right)\)
    priorityList
    \(\sigma=\frac{e_{\max }-e_{\text {min }}}{2}\)
    foreach \(A_{i} \in \mathcal{M}^{\prime}\) do
        if \(\left|A_{i}\right|>m_{i}+\sigma\) then
        priorityList.Add \(\left(A_{i}\right)\)
    else if \(\left|A_{i}\right|<m_{i}-\sigma\) then
                priorityList.Add \(\left(A_{i}\right)\)
    end if
    end foreach
    SortDescent (priorityList)
    return priorityList
```

traverses the list verifying the type of each edge and applies the appropriate operation (edge split or edge collapse). All the vertices adjacent to an edge that was changed are marked as visited, if both vertices of the edge were visited than the edge is removed from the list $L_{p}$ (Algorithm 3). This imposes a even more smoothly change between $\mathcal{M}_{i}^{\prime}$ and $\mathcal{M}_{i+1}^{\prime}$. Note that an edge can be in the list $L_{p}$ and after an operation that affect one of its vertices not be long or short .

```
Algoritmo 3: StellarOperations \(\left(\mathcal{M ~}^{\prime}, L_{p}\right)\)
    foreach \(A_{i} \in L_{p}\) do
        if BothVertexVisited \(\left(A_{i}\right)\) then
            continue
        else if \(\left|A_{i}\right|>m_{i}+\sigma\) then
            EdgeSplit \(\left(A_{i}\right)\)
            VisitEdgeNeighbors \(\left(A_{i}\right)\)
        end if
        else if \(\left|A_{i}\right|<m_{i}-\sigma\) then
            EdgeCollapse ( \(A_{i}\) )
            VisitEdgeNeighbors \(\left(A_{i}\right)\)
        end if
    end foreach
```

In this work we developed a method to reposition the output vertices of edge split and edge collapse operations. It position the vertex over the original edge in a way that
minimize the equalization:

$$
\begin{equation*}
\sum_{V_{j}} A B S\left(\left|V_{i}-V_{j}\right|-m_{i}\right)^{2} \tag{3.1}
\end{equation*}
$$

where $V_{i}$ is the vertex we want to position and $V_{j}$ are the vertices connected to $V_{i}$.

### 3.2 Valency optimizer

After the edge adjustments by stellar operations, the valency of the vertices naturally tends to 6 , which is mandatory to obtain uniform distribution in smooth regions. However, some vertices might have arbitrary valencies. A simple scheme to correct the valencies of the vertices based on edge flip operations is then applied (Mario et al, 2010a).

Consider the two triangles of Figure 2.2. For each edge of the mesh, the algorithm performs an edge flip operation and, if the valencies of the involved vertices become near to 6 and the angle $\min \left(\alpha_{i}\right)$ is greater than $f \cdot \min \left(\alpha_{i}\right)$ before the edge flip, the rotated edge is accepted, otherwise the operation is undone. An overview of the process is shown on Algorithm 4.

```
Algoritmo 4: CorrectValency ( \(\mathcal{M}^{\prime}\) )
    foreach \(A_{i} \in \mathcal{M}^{\prime}\) do
        \(\operatorname{preMin}_{\alpha}=\operatorname{getMinalpha}\left(A_{i}\right)\)
        \(v_{n}=\) SearchAdjacentVertices \(\left(a_{i}\right)\)
        preDeviation \(=\left|\operatorname{valency}\left(V_{1}\right)-6\right|+\mid\) valency \(\left(V_{2}\right)-6\left|+\left|\operatorname{valency}\left(V_{3}\right)-6\right|\right.\)
        \(+\mid\) valency \(\left(V_{4}\right)-6 \mid\)
        EdgeFlip \(\left(A_{i}\right)\)
        postMin \({ }_{\alpha}=\) getMinalpha \(\left(A_{i}\right)\)
        postDeviation \(=\mid\) valency \(\left(V_{1}\right)-6|+|\) valency \(\left(V_{2}\right)-6 \mid+\)
        \(\mid\) valency \(\left(V_{3}\right)-6|+|\) valency \(\left(V_{4}\right)-6 \mid\)
        if postDeviation \(\leq\) preDeviation and postMin \({ }_{\alpha} \leq f\).preMin \({ }_{\alpha}\) then
            EdgeFlip \(\left(A_{i}\right)\)
        end if
    end foreach
```


### 3.3 Lowpass Filtering

After the valency optimizer step, we proceed to the lowpass filtering. In this step, we use an iterative method to approach the constrained Laplacian. In the classical Laplacian filter, we add some additional constraints to reduce the geometry loss:

$$
\begin{array}{r}
N_{i} \cdot T_{i}=0, \forall T_{i} \in \mathcal{M}^{\prime}, \\
\left|T_{i}\right|=0 \forall T_{i} \in \mathcal{B},
\end{array}
$$

where $N_{i}$ is the normal of the current mesh in the vertex $V_{i}$ and $T_{i}$ is the unknow displacement of the vertex $V_{i}$.

```
Algoritmo 5: LowPassFiltering \(\left(\mathcal{M}^{\prime}, k\right)\)
    foreach \(V_{i} \in \mathcal{M}^{\prime}\) do
        \(k S t a r=\operatorname{getKStar}\left(V_{i}, k\right)\)
        fat \(=0\)
        foreach \(V_{j} \in k S t a r\) do
            \(V_{i}^{\prime}+=\frac{V_{j}}{s t a r}\)
            fat \(+=\frac{1}{s t a r}\)
        end foreach
        \(V_{i}^{\prime}=\frac{V_{i}^{\prime}}{f a t}\)
        \(T_{i}=V_{i}^{\prime}-V_{i}\)
        \(T_{i}=T_{i}\)-projection \(\left(T_{i}, N_{i}\right)\)
    end foreach
    foreach \(V_{i} \in \mathcal{N}^{\prime}\) do
        if \(V_{i} \notin \mathcal{B}\) then
            \(V_{i}+=T_{i}\)
        end if
    end foreach
```

The iterative Algorithm 5 approximate the constrained Laplacian filtering described above. It calculates the new vertex position based on the $k$-neighborhood as proposed in Oliveira et al (2012). The first step is to compute for each vertex the new position without the application of the new constraints. This position is defined by the center of mass of all neighbors vertices weighted by their ring number in such a way that distant vertices contribute less than near vertices.

The second step is to impose the constraint $N_{i} \cdot T_{i}=0$ by removing the vector
corresponding to the projection of $T_{i}$ in $N_{i}$. When all displacements are computed, the vertices $V_{i}$ are updated except those on the borders.

### 3.4 Projection

After the previous steps the new vertices tend to move away from the original mesh. As proposed in (Oliveira et al, 2013), this problem can be reduced by projecting the mesh computed in each iteration over the original mesh. However, the projection of a mesh over another is a complicate problem since the object can have arbitrary topology and geometry. Even if the mesh to be projected is close to the original, artifacts may appear due to wrong vertex correspondences. The algorithm works as follows: for each vertex we compute its nearest projection over some original triangles. These triangles are the incident over the 30 nearest vertices in the original mesh (Algorithm 6). To justify the 30 nearest vertices, we run several tests where using 30 nearest neighbors the resulting projection was exactly the same as if we project in all triangles. But this behaviour is not guaranteed for all cases because it is dependent on the geometry and topology of the mesh.

```
Algoritmo 6: Projection( \(\left.\mathcal{M}, \mathcal{M}^{\prime}\right)\)
    foreach \(V_{i} \in \mathcal{M}^{\prime}\) do
        Nearest \(=\) getNearest \(\left(V_{i}, \mathcal{M}, 30\right)\)
        foreach Triangle \(_{j} \in\) Nearest do
            \(P=\) projectionVT \(\left(V_{i}\right.\), Triangle \(\left._{j}\right)\)
            if \(\left|V_{i}-P\right| \leq\left|V_{i}-M I N\right|\) then
                \(M I N=P\)
            end if
        end foreach
        \(V_{i}=M I N\)
    end foreach
```

In our implementation of the Algorithm 6, we use a Kd-tree to increase the performance.

## 4 Results

In this section, the results of this work are shown. The method was implemented using C++ programming language and compiled using GCC 4.6.3. All tests were run in a Intel Xeon(R) CPU E31220 @ $3.10 \mathrm{GHz} \times 4$ computer that has 8 GBs of RAM. The graphic card was an AMD Radeon HD 5700 series.

### 4.1 Edge equalization by simplification

In this first scenario, we were testing the iterative method evolution in time (Table 4.1). The following tests were run over the fertility model (Figure 4.1). The parameters used in the tests were $n=100, k=2, e_{\text {min }}=1.2, e_{\text {min }}=1.8, p r=0, f=\frac{1}{2}$ and the input mesh has an original edge length average of 0.75.


Figura 4.1: This is the input fertility model.

The table 4.1 columns are the iteration Iteration, the number of vertices Vertices,

| Iteration | Vertices | Edges | Average | Deviation | Regular | Small | Big |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 9354 | 28080 | 0.934116 | 0.150386 | 72.011973 | 26702 | 3 |
| 6 | 3629 | 10905 | 1.488004 | 0.199443 | 73.215762 | 889 | 505 |
| 11 | 3492 | 10494 | 1.513421 | 0.187043 | 77.348225 | 623 | 477 |
| 16 | 3464 | 10410 | 1.518211 | 0.171593 | 79.907621 | 458 | 339 |
| 21 | 3425 | 10293 | 1.525191 | 0.163202 | 80.350365 | 369 | 304 |
| 26 | 3401 | 10221 | 1.529626 | 0.161322 | 81.770068 | 360 | 284 |
| 31 | 3400 | 10218 | 1.530237 | 0.165183 | 80.970588 | 377 | 291 |
| 36 | 3384 | 10170 | 1.532931 | 0.163540 | 82.092199 | 381 | 321 |
| 41 | 3405 | 10233 | 1.528674 | 0.159297 | 82.085169 | 362 | 274 |
| 46 | 3394 | 10200 | 1.530222 | 0.161740 | 81.644078 | 382 | 299 |
| 51 | 3382 | 10164 | 1.532950 | 0.157780 | 82.820816 | 340 | 277 |
| 56 | 3361 | 10101 | 1.536691 | 0.150204 | 84.141625 | 272 | 238 |
| 61 | 3380 | 10158 | 1.533097 | 0.151148 | 84.171598 | 300 | 228 |
| 66 | 3381 | 10161 | 1.532640 | 0.150045 | 85.241053 | 267 | 267 |
| 71 | 3382 | 10164 | 1.532510 | 0.151162 | 83.973980 | 300 | 217 |
| 76 | 3376 | 10146 | 1.533919 | 0.152444 | 84.360190 | 299 | 241 |
| 81 | 3371 | 10131 | 1.535301 | 0.148051 | 84.307327 | 270 | 240 |
| 86 | 3391 | 10191 | 1.530824 | 0.151205 | 84.370392 | 304 | 240 |
| 91 | 3416 | 10266 | 1.525567 | 0.159351 | 83.313817 | 393 | 253 |
| 96 | 3375 | 10143 | 1.534117 | 0.148669 | 85.214815 | 276 | 237 |

Tabela 4.1: This table shows the iterative method evolution in time. It clearly shows the refinement in the mesh, with the decreasing number of vertices.
number of Edges, the average edges length Average, the standard deviation of the average Average, the percent of regular vertices Regular, the number of edges that are smaller than the interval Small and the number of edges that are bigger than the interval Big.

As depicted in Figure 4.2 the method generates a almost uniform mesh without losing geometry accuracy. Even the trivalent mesh is close to the original geometry and almost uniform as depicted in Figure 4.3.

### 4.2 Edge equalization by refinement

As edge equalization by simplification we run tests for measuring the iterative method evolution in time (Table 4.2). To be more accuracy, the following results were running in the same fertility model used in edge equalization by simplification (Figure 4.1).The parameters used in the tests were $n=100, k=2, e_{\text {min }}=1.2, e_{\text {min }}=1.8, p r=0, f=\frac{1}{2}$ and the input mesh has an original edge length average of 2.25 .

As depicted in Figure 4.4 the method generates a almost uniform mesh without


Figura 4.2: The first mesh is the input mesh, the second one is the mesh in the 100th iteration.


Figura 4.3: The first mesh is the input mesh, the second one is the hexagonal mesh in the 100th iteration.
losing geometry accuracy. As depicted in Figure 4.5 even the trivalent mesh is close to the original geometry and almost uniform.

| Iteration | Vertices | Edges | Average | Deviation | Regular | Small | Big |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | 17539 | 52635 | 2.045783 | 0.327601 | 67.478191 | 394 | 40852 |
| 6 | 32826 | 98496 | 1.492026 | 0.195853 | 72.156218 | 7353 | 4301 |
| 11 | 31479 | 94455 | 1.518870 | 0.164277 | 78.843038 | 3833 | 2717 |
| 16 | 31058 | 93192 | 1.526876 | 0.147635 | 82.326615 | 2558 | 1725 |
| 21 | 30690 | 92088 | 1.534388 | 0.133350 | 84.874552 | 1731 | 1154 |
| 26 | 30491 | 91491 | 1.538608 | 0.124447 | 86.199206 | 1225 | 729 |
| 31 | 30374 | 91140 | 1.541098 | 0.119997 | 87.041549 | 991 | 584 |
| 36 | 30289 | 90885 | 1.542784 | 0.115151 | 87.817359 | 740 | 420 |
| 41 | 30236 | 90726 | 1.543908 | 0.113076 | 88.176346 | 634 | 357 |
| 46 | 30195 | 90603 | 1.544772 | 0.110984 | 88.491472 | 522 | 363 |
| 51 | 30192 | 90594 | 1.544682 | 0.109332 | 88.665872 | 491 | 232 |
| 56 | 30124 | 90390 | 1.546264 | 0.107284 | 88.965609 | 351 | 224 |
| 61 | 30129 | 90405 | 1.546056 | 0.106334 | 89.146669 | 326 | 183 |
| 66 | 30114 | 90360 | 1.546384 | 0.106494 | 89.144584 | 328 | 188 |
| 71 | 30080 | 90258 | 1.547069 | 0.104416 | 89.507979 | 238 | 125 |
| 76 | 30076 | 90246 | 1.547142 | 0.104279 | 89.559782 | 214 | 129 |
| 81 | 30085 | 90273 | 1.546890 | 0.103804 | 89.566229 | 208 | 103 |
| 86 | 30083 | 90267 | 1.546922 | 0.103874 | 89.572184 | 209 | 97 |
| 91 | 30068 | 90222 | 1.547238 | 0.103024 | 89.696688 | 159 | 102 |
| 96 | 30084 | 90270 | 1.546902 | 0.103884 | 89.612419 | 204 | 109 |

Tabela 4.2: This table shows the iterative method evolution in time. It clearly shows the refinement in the mesh, with the increasing number of vertices.


Figura 4.4: The first mesh is the input mesh, the second one is the mesh in the 100th iteration.


Figura 4.5: The first mesh is the input mesh, the second one is the hexagonal mesh in the 100th iteration.


Figura 4.6: The first mesh is a simplified output mesh, the second one refined output mesh.

The edge equalization by refinement is better than by simplification because it preserve more the original geometry as depicted in Figure(4.6). Due this, the graphics
were based on edge equalization by refinement.

### 4.3 Per iteration graphs

The graphics in this section were generated using the following parameters $n=100, k=$ $2, e_{\text {min }}=1.2, e_{\text {min }}=1.8, p r=0, f=\frac{1}{2}$ and the input mesh has an original edge length average of 2.25 .


Figura 4.7: The boxes represent the time spent in each iteration.

As this is a refinement the number of edges and vertices increases over the early iterations, so the time spent increases as well. After a peak the time spent per iteration stabilizes around nine seconds as depicted in Figure 4.7.

The edge length average quickly converge to a value next to $m=\frac{e_{\min }+e_{\max }}{2}$ as depicted in Figure 4.8. The method changes the mesh so fast that the average oscillates around the expected average $m$ before converging. Due to these drastic changes, the standard deviation decreases slowly to its minimum.

As depicted in Figure 4.9, the number o edges lengths out of interval decreases even more slowly than the standard deviation. this was expected as the imposition of a minimum and maximum value for edge lengths is much more restrictive since the majority of the edges cannot compensate the extra size of a few outliers.


Figura 4.8: The first curve shows the edge length average in per iteration. The second curve is the standard deviation.

In Figure 4.10, we can see the method adapting the number of vertices in the model according to the constraints.

In Figure 4.11, we can see that the valency optimizer really makes a good work as the number of regular vertices slowly rises until it is more than $90 \%$.

As depicted in Figure 4.12, the dual mesh is almost an uniform hexagonal mesh. There are almost no edges lengths out of the predefined interval. This is very important due the applications of this kind of mesh in physics.


Figura 4.9: The red curve shows the number of edges lengths out of the interval per iteration. The green curve shows the number of small edges. The blue curve shows the number of big edges.


Figura 4.10: This shows the number of vertices per iteration.

### 4.4 Flaw case

This method can not do a strong simplification, as this will remove so much vertices, in such a way that it can not sample the entire surface. In the example (Figure 4.13) the average edges lengths were five times smaller than the average of the predefined interval.


Figura 4.11: This shows the percent of regular vertices per iteration.


Figura 4.12: This is the histogram of the hexagonal mesh edges length at 100th iteration.

### 4.5 Examples

These are other examples of meshes processed using this method. The tests were run over the Rockarm (Figure 4.14) and Bunny (Figures 4.17,4.15,4.18) models with the same parameters $n=50, k=2, e_{\text {min }}=1.2, e_{\text {min }}=1.8, p r=0, f=\frac{1}{2}$ and the input mesh has an original edge length average of 2.25 .

The method is robust and can be applied even on larges meshes with thousands of vertices. The Rockarm model (Figure 4.14) has 20776 vertices and after the process


Figura 4.13: This is a flaw case of the method. The output mesh on the bottom loses almost the entire original mesh geometry.


Figura 4.14: On the top is the input mesh, on the bottom the mesh after 50 iterations.
the output hexagonal model has 95020 vertices.
The method is powerful enough to handle extremely irregular meshes, as in the


Figura 4.15: On the top is the input mesh, on the bottom the mesh after 50 iterations.


Figura 4.16: The same model depicted in Figure 4.15, with the ear is in focus.
bunny model depicted in Figures 4.15, 4.16, 4.17 and 4.18, this model has higher frequency regions as depicted in Figure 4.16 and vertices with extremely high valencies and edges with lengths much bigger or much smaller than the average as depicted in Figure 4.17.


Figura 4.17: On the top is the input mesh, on the bottom the mesh after 50 iterations.


Figura 4.18: This is the trivalent bunny after 50 iterations.

## 5 Conclusion

This work proposes a low memory cost edge length equalization method. The edge length is an important mesh quality for physics simulations. Due to this, a method capable of edge equalization may be very useful as there are many publicly models that could be used in simulations after processed.

The method will not converge if the predefined interval is too tide. In our tests a good interval have to be enough spaced in such a way that the stellar operations will be applied rarely in the later iterations. This is often an interval obtained by fixing an average $m$, setting $e_{\text {min }}=m-0.2 m$ and setting $e_{\text {min }}=m+0.2 m$.

When the algorithm diverges, it is possible to see that the number of vertices in the priority list will continue high over the iterations and the projection will move vertices too far from their original positions. In that case setting a larger interval might reduce or solve this problem. Another problem is when the target interval forces the model to be extremely simplified.

Besides the $e_{\min }$ and $e_{\max }$, the method is very sensitive to some parameters. The $k$ rings can be fixed at $k=2$, as our tests suggests that higher $k$ does not improve very much the algorithm but increases the computational effort.

The frequency of projection per iteration $p r$ is very dependent of the desired objective because it allows a greater geometry loss, but helps the method to converge in a more tide interval.

The maximum angle for flipping operations $f$ does not affect the convergence very much, and a value of $f=0.5$ gives good results in most cases. This prevents a great geometry loss but allows the algorithm to flip edges for improving the vertex valency in a balanced way.

### 5.1 Future works

This method needs a post process to remove few edges that are not within the predefined interval. Another thing to do is the implementation of another lowpass filter, for comparison with the actual Laplacian filter or even design a combined filter.

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