# Optimal Good-Aspect-Ratio Coarsening for Unstructured Meshes

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#### Abstract

A hierarchical gradient of an unstructured mesh  $M_0$ is a sequence of meshes  $M_1, \ldots, M_k$  such that  $|M_k|$ is smaller than a given threshold mesh size b. The gradient is well-conditioned if for each i in the range  $1 \leq i \leq k$ , (1)  $M_i$  is well-shaped, namely, elements of  $M_i$  have a bounded aspect ratio; and (2)  $M_i$  is a coarsened approximation of  $M_{i-1}$ . The gradient is node-nested if the set of the nodes of  $M_i$  is a subset of that of  $M_{i-1}$ . The problem of constructing well-conditioned coarsening gradients is a key step for hierarchical and multi-level numerical calculations. In this paper, we give an algorithm for finding a wellconditioned hierarchical gradient of a two dimensional unstructured mesh. Our algorithm can be used to generate both node-nested and non-nested gradients. The gradient  $M_1, \ldots, M_k$  we generate is optimal in the following sense: there exists a constant c such that for any other well-conditioned hierarchical gradient  $M'_1, \ldots, M'_k, |M_i| \leq c|M'_i|$ , that is, the size of the mesh at each level is smaller up to a constant factor.

### 1 Introduction

The class of hierarchical and multi-level techniques has become one of the most effective and successful classes of numerical techniques for solving partial differential equations (PDEs). These techniques have been used in multigrid methods [4] and multi-level domain decomposition [5].

Numerical methods such as the finite element, finite difference, and finite volume methods apply the following five basic steps to solve a PDE over a domain  $\Omega$ .

- Formulate the problem (e.g., in term of PDEs) and describe the geometry and the boundary of a continuous domain D (sometime called geometric modelling).
- 2. Generate a well-shaped mesh M to approximate

the domain.

- 3. Generate a system of linear or non-linear equations over M for the governing PDEs (e.g., assemble the stiffness matrix and the right hand vector).
- 4. Solve the system of equations and estimate the error of the solution.
- 5. Adaptively refine the mesh and return to step (3) if needed.

Once the mesh M is generated, we need to solve a system of linear equations defined over M. A hierarchical method solves this linear system by first constructing a hierarchical gradient of meshes  $M_0, \ldots, M_k$ , where  $M_0 = M$  is the finest mesh that discretizes  $\Omega$ . For each i in the range  $1 \leq i \leq k$ , the mesh  $M_i$  is a geometric coarsening of  $M_{i-1}$ . The gist of multigrid methods and other hierarchical numerical methods is the transformation of partial solutions from mesh  $M_i$  to mesh  $M_{i-1}$ using interpolation, and from mesh  $M_i$  to mesh  $M_{i+1}$ using restriction. Informally, these hierarchical methods solve a PDE on  $\Omega$  by first obtaining an initial vector solution either for  $M_0$  or for  $M_k$ , and then improving the quality of the vector by transforming it hierarchically up and down the hierarchy while applying some simple and efficient iterative methods at each level.

The simplest form of a hierarchical gradient is a series of nested structured meshes (regular grids). Brandt showed, by carefully using restriction and interpolation, that the solution for  $M_0$  can be obtained very efficiently using multigrid methods. See also Bramble, Pasciak and Xu [10]. Nested structured hierarchical gradients are attractive choices in practice because they can be easily generated and because the convergence of the structured multigrid methods is well understood. However, the use of structured regular grids limits the applicability of this simplest class of hierarchical gradients to problems whose domains are simple and whose solution functions have small or constant Hessian [12, 5, 18].

The use of unstructured meshes is inevitable in the solution of complex problems with more intricate domain geometry and solutions. This paper concerns with the problem of generating quality hierarchical gradients for unstructured meshes.

The effectiveness of a hierarchical method that uses

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an unstructured hierarchical gradient  $M_0,\ldots,M_k$  depends on the quality of this gradient [10, 5, 6]. In particular, Chan and Zou provided sufficient conditions for multilevel additive Schwarz methods to work on unstructured meshes. Informally, their conditions require, for each i in the range  $1 \leq i \leq k$ , that (1)  $M_i$  is well-shaped, e.g., in two dimensions elements of  $M_i$  should have a bounded aspect ratio; and (2)  $M_i$  approximates  $M_{i-1}$  in the numerical formulation. The coarsening problem can thus be informally defined as: Given a well-shaped mesh  $M_0$  and a threshold size b, construct a gradient  $M_1,\ldots M_k$  with  $|M_k| \leq b$  that satisfies conditions (1) and (2).

In this paper, we give an algorithm for the coarsening problem. Our algorithm guarantees the quality of the hierarchical gradient (i.e., it produces gradients that satisfies conditions (1) and (2)). It also minimizes the size of the mesh at each level up to a constant factor. A gradient is node-nested if the set of the nodes of  $M_i$  is a subset of that of  $M_{i-1}$ . Our algorithm can be used to generate both node-nested and non-nested coarsening gradients.

A version of this research, for the simpler case of quasi-uniform mesh coarsening, was included in a survey of our work we submitted to the 5th international meshing roundtable [13]. In this paper we address the problem of general unstructured mesh coarsening.

### 2 The problem of Mesh Coarsening

2.1 Mesh qualities. A two dimensional domain  $\Omega$  is a planar straight-line graph (PSLG), whose boundary is polygonal. The edges in the PSLG can represent boundaries between two materials, points of special interest or holes. In this extended abstract we assume a simple form of this general definition. Our domain is the unit square, and its boundary is the square's four edges.

A mesh is a discretization of the domain. The discrete components of the mesh are the mesh elements. We mostly discuss triangular meshes, whose elements are triangles. We will often refer to the following three categories of meshes: (1) grids: a mesh whose elements are squares of equal size. (2) quasi-uniform unstructured meshes: a mesh whose elements' side lengths differ by at most a constant factor. (3) unstructured meshes: a general mesh, with no restriction of the elements' size and shape.

Not all meshes perform equally well in numerical computations. Numerical and discretization error depend on the geometric shape and size of the mesh elements [17]. We will use the following definitions for quantifying the geometric shape and mesh element size.

Definition 2.1. The edge-length function of a

 $\label{eq:mesh} \begin{array}{l} \textit{mesh } M, \ \textit{el}_{M}, \ \textit{is defined for each} \ x \in \Omega \ \textit{to be the length} \\ \textit{of the longest edge of all the mesh elements that contain} \\ \\ \textbf{x} \end{array}$ 

DEFINITION 2.2. The aspect ratio of a triangular mesh element is the smallest angle of the element. The aspect ratio of a mesh is the smallest angle of its elements.

A mesh is said to be of bounded aspect ratio if its aspect ratio is larger than  $\theta$ , where  $\theta$  is a predefined parameter quantifying the mesh quality. We note that there are many definitions for the aspect ratio of a two dimensional mesh which are interchangeable with the above definition [2].

The *el* function is a measure of the mesh element size; the aspect ratio is a measure of the element shape. These two qualities are often at odds: to produce a good aspect ratio mesh conforming to the boundary the mesh often has to be finer. The goal of *mesh generation* is to construct a good aspect ratio mesh whose element sizes are as large as possible (so that the number of mesh elements is as small as possible).

**2.2** Mesh coarsenings. A coarsening M' of a mesh M is a mesh whose edge length function  $el_{M'}$  is pointwise bigger than  $el_M$  but still conforms to the same domain

The coarsening can be classified as element-nested, node-nested or non-nested. In general, a triangular mesh does not have any element-nested coarsening, unless it was carefully crafted as such. Furthermore, coarsening, even in the relaxed sense of node-nested meshes, can cause a degradation in the aspect ratio of the coarser mesh compared to that of the finer mesh. In this paper, we are interested in generating a sequence of respectively coarser meshes - a gradient of mesh coarsenings. This sequence will be referred to as the  $hierarchical\ coarsening\ gradient$  of M. The objective of this paper is to develop an automatic mesh coarsener which guarantees good aspect ratio of the entire coarsening gradient. We now introduce some definitions to formalize our discussion.

DEFINITION 2.3. The **depth** of an hierarchical gradient is the number of meshes (levels) in the gradient; The **width** of level i of the gradient is  $|M_i|$ , the number of elements in the mesh  $M_i$ .

DEFINITION 2.4. Let  $\theta$  be a constant in the range  $0 < \theta < \pi/2$ ; and let b, I be two positive constants. A hierarchical gradient  $M_0, ..., M_k$  in two dimensions is a  $(\theta, \mathcal{I}, b)$ -Well-Conditioned Hierarchical gradient if  $|M_k| \leq b$  and for each i in the range  $1 \leq i \leq k$  each angle  $\beta$  in  $M_i$  satisfies  $\theta < \beta < \pi - \theta$ . Furthermore, every two adjacent meshes are  $\mathcal{I}$ -locally similar, i.e.  $el_{M_{i+1}} \leq \mathcal{I}el_{M_i}$ .

The first condition, good aspect ratio, is motivated by the requirement from iterative methods: The multigrid method and multi-level domain decomposition use an iterative method to "smooth" the residual error at each level. The convergence properties of iterative methods are related to the aspect ratio of the underlying mesh. Chan and Zou [6] showed that bounded aspect ratios at each level are important for both additive Schwarz based multi-level domain decomposition and multigrids. The second condition, the local similarity, is motivated by the restriction and interpolation phases of the multigrid methods, which are used to transform partial solution between meshes in adjacent levels of the hierarchy. To reduce the interpolation and restriction errors, adjacent meshes should approximate each other well. Both local similarity and bounded aspect ratio were used in Chan and Zou's analysis, which showed that they are sufficient for multilevel additive Schwarz methods to work on unstructured meshes.

Given  $M_0$  and  $(\theta, \mathcal{I}, b)$ , the problem of **hierarchical** mesh coarsening is to find a  $(\theta, \mathcal{I}, b)$ -well conditioned hierarchical gradient with smallest depth and width.

### 2.3 Previous approaches to mesh coarsening.

To produce a node nested coarsening, a subset of the fine mesh nodes is picked, and retriangulated to form the coarse mesh. Various approaches to mesh coarsening differ in the methods they use to pick the coarser mesh nodes, and the retriangulation method. The problem of mesh coarsening received much attention, see [9, 19, 11, 15, 1, 7]. However, none of the papers we found address the geometrical issues of mesh coarsening. Properties such as element quality of the mesh or the size of the intersection of the coarse and fine mesh (their local similarity) are not discussed beyond empirical observations.

Picking the node set of the coarser mesh using a maximal independent set (MIS) technique seems to be the most popular approach [5, 9, 19, 11, 15, 7]. The 1D skeleton of the fine mesh is viewed as a graph, and a set of nodes such that no two share an edge is picked (independence). A node not picked for the coarse mesh must neighbour a node that was retained, so the independent set is maximal. An MIS can be constructed to ensure a constant factor reduction in the number of the mesh nodes. However, we now show that the MIS technique can not guarantee the aspect ratio of the coarsened meshes, both for quasi-uniform and for general unstructured meshes.

Quasi-uniform unstructured mesh coarsening: The MIS on the 1D skeleton technique is very successful in reducing the mesh size to a fraction of its original size; however it carries no guarantees for the other

qualities of the mesh hierarchy, such as its aspect-ratio. The problem is illustrated in Figure 1: certain choices of an MIS of the original mesh degrade the aspect ratio of the coarser mesh. The aspect ratio degradation compounds with repeated applications. This can be observed even for very uniform meshes, as in the grid-like mesh of the figure.







Figure 1: Repeated applications of MIS can degrade the aspect ratio.

# General unstructured mesh coarsening:

Above we gave an example of a quasi-uniform mesh for which certain choices of the MIS cause repeated degradation of its aspect ratio. For general unstructured meshes, a much stronger statement is true: there exists an unstructured mesh such that all possible choices of MIS result in a coarsening hierarchy with increasingly worse aspect ratio.

Let  $M_0$  be a one dimensional mesh whose nodes are  $P = \{2^{i-1} - 1 : i = 1 \cdots n\}$ . The edges of the mesh are between adjacent points,  $M_0$  is a line graph of n nodes. The aspect ratio of a one dimensional mesh is the maximum ratio between two adjacent edges, hence the aspect ratio of  $M_0$  is equal to 2.

For this mesh there is no well-conditioned coarsening gradient of depth  $\log n$ . This statement is a corollary of the results presented in this paper. Since the MIS technique reduces the size of the mesh by a constant factor at each level, it fails to produce good aspect ratio coarsenings.

In particular, the MIS technique defines a hierarchical gradient where the jth mesh  $\{M_j\}$  has node set  $P_j = \{2^{i2^j-1} - 1 : i = 1 \cdots n/2^j\}$ . In other words,  $P_j$  is formed by taking every other point of  $\mathcal{P}_{j-1}$ . The aspect ratio of mesh  $M_j$  is therefore at least  $2^{2^j} - 1$ , and worsens super–exponentially. This one dimensional example can easily be extended to two and three dimensions.

As part of the results in this paper, we will obtain bounds for the shortest (up to a constant) well-conditioned gradient for this mesh and provide simple algorithms for its generation.

### 3 Function Based Mesh Coarsening

In this section, we present our approach to mesh coarsening. In order to construct a well-conditioned hierarchical gradient we have to overcome the escalating degradation of the mesh quality, demonstrated in the previous section.

Rather than using only the information present in the respectively coarser meshes, we use an intermediate representation which we term "spacing functions". The spacing functions, along with the point set of the initial mesh, capture the information necessary for the generation of the coarser meshes.

At a high level,  $M_0$  defines a spacing function  $f_0$  which describes the typical size and point spacing of  $M_0$ . Our idea is to compute a spacing function  $f_i$  for each level, and use it to generate  $M_i$ . Given the spacing functions our task is then to create a point set that is "spaced" according to that function, and triangulate it (as for mesh generation [12]). We refer this mesh coarsening technique as **function-based coarsening**. It contains four steps: (1) recover the spacing function of the initial mesh; (2) increase the spacing value of the mesh points smoothly to obtain the new spacing functions; (3) delete some mesh nodes so that the remainder nodes are spaced according to the new spacing function; and (4) compute the Delaunay triangulation of the nodes obtained in Step (3).

# **3.1 Recovering the spacing function:** We first formalize our notion of a spacing—function.

DEFINITION 3.1. Let  $\beta > 1$  be a real number. A point set P is  $\beta$ -spaced according to a function f if for any two points  $p, q \in P$ ,  $f(p_1) + f(p_2) < \beta ||p - q||$ . The function f is then referred to as the  $\beta$ -spacing function of P.

The initial spacing function we use is based on the natural spacing of the original mesh:

DEFINITION 3.2. The nearest neighbour (NN) function of a point set  $P \subset \Omega$  assigns to each point  $p \in P$  the distance to the point  $q \in P$  nearest to it such that  $q \neq p$ . It can be extended to the domain  $\Omega$  by assigning to a point  $x \in \Omega$  the radius of the smallest closed ball centered at x and containing at least two points from P.

### 3.2 Coarsening the spacing functions:

DEFINITION 3.3. Let P be a point set in a domain  $\Omega$  in  $\mathbb{R}^2$ . Let g be a spacing function over  $\Omega$ . Let C > 1 be a real number. The C-coarsening of f with respect to P is a spacing function over  $\Omega$  such that for all  $x \in \Omega$ 

$$f_{g,C,P}(x) = \min_{p \in P} \hat{f}_{g,C,p}(x),$$

where for each point  $p \in P$ ,  $\hat{f}_{g,C,p}(x) = C \cdot g(p) + \|p-x\|$ . When clear from the context, we omit g from the notation.

We can use a simpler way to generate the coarsening function for quasi-uniform functions g:

**Procedure:** ONE\_LEVEL\_COARSEN(M)

Input: M, a mesh over a square  $\Omega$ . C, the coarseing factor.

g, the  $\beta$ -spacing function of M.

Output:  $M_1$ , a coarser triangular mesh. Method:

- 1. Let  $P_0$  be the square's corners;  $P_1$  the mesh nodes located on the square's edges;  $P_2$  the rest of the mesh nodes.
- 2. Compute  $f(p_i) = f_{g,C,P}(p_i)$  for each  $p_i \in P$ .
- 3. Let  $\hat{P}_2$  be the set of points  $\{p\}$  in  $P_2$  whose distance from the boundaries is at least  $\delta f(p)$ . ( $\delta$  is a small fixed constant.)
- 4. Construct a conflict graph with respect to  $f\colon \ CG(P_0\cup P_1\cup \hat{P_2})$ .
- 5. Let S be a maximal independent set of CG generated by first considering the points  $P_0$ , then  $P_1$ , and finally  $\hat{P}_2$ .
- 6. Return  $M_1 \, = \, DT(S) \, , \ \, \mbox{the Delaunay triangulation} \label{eq:mass_state}$  of  $S \, .$

Figure 2: One level function based coarsening.

DEFINITION 3.4. Let P be a point set in a domain  $\Omega$  in  $\mathbb{R}^2$ . Let g be a spacing function over  $\Omega$ . Let C>1 be a real number. The C-threshold-coarsening function is defined as:

$$t_{g,C,P}(x) = \max(g(x), C \min_{y \in \Omega} g(y))$$

The set of coarsening functions we suggest is therefore:

- for general coarsening:  $\{f_{NN,2^i,P}\}$  and  $\beta$  is a constant depending on the aspect ratio (see section 5).
- for quasi–uniform coarsening:  $\{t_{NN,2^i,P}\}$  and  $\beta=2.$
- 3.3 Coarsening the meshes: Let  $M_0$  be the initial mesh, C the factor by which the mesh should be coarsened. To coarsen the mesh, we first generate the coarsened spacing function values for each node of the mesh using C, and then pick a subset of the mesh nodes which is  $\beta$ -spaced by this spacing function. One possible method to pick the coarsened mesh nodes is by using a conflict graph.

DEFINITION 3.5. The **conflict graph** of a point set P, CG(P), with respect to a  $\beta$ -spacing function f is a graph CG(P) = (P, E) where

$$E = \left\{ (p_i, p_j) : ||p_i - p_j|| < \frac{f(p_i) + f(p_j)}{\beta} \right\}$$

Our proposed scheme is outlined in Figure 2. This one-level scheme can be naively extended to a multi-level scheme by repeatedly applying it with coarser and coarser spacing functions  $f_i$ , generated using function coarsening and constants of the form  $C^i$  for level i.

However this in general produces a gradient that is not node-nested, for it generates the mesh at each level independently of other levels. To construct a node-nested hierarchical gradient, we need a subtler approach. At a high level, we first generate  $M_k$ , the coarsest mesh. We then enforce that nodes in  $M_k$  will be chosen in  $M_{k-1}$ . Repeating this enforcement, we can build a node-nested well-conditioned gradient. The resulting scheme is outlined in Figure 3.

Lemma 3.6.  $S_{i+1}$  is also an independent set of  $CG_i$ .

**Proof**: Because C > 1,  $C^{i+1} > C^i$ . Therefore, for each point  $x \in \Omega$ ,  $f_{C^{i+1},P}(x) \ge f_{C^i,P}(x)$ . By definition 3.5,  $CG_{i+1}$  is a supergraph of  $CG_i$ . Hence any independent set of the former must be independent in the later as well.

Therefore, Algorithm MULTI\_LEVEL\_COARSEN correctly generates a node-nested multilevel gradient.

The rest of this paper will show the correctness of this coarsening method, stated in the following two theorems:

THEOREM 3.7. Let  $M_0$  be a mesh whose smallest angle is bounded bellow by  $\theta$ . The hierarchical gradient  $(M_1, ..., M_k)$  produced by the algorithm of Figure 3 has the following properties:

- 1. **aspect ratio:** There is a constant  $\theta_1$  depending on  $\theta$  only such that for  $1 \leq i \leq k$ , the smallest angle of mesh  $M_i$  is bounded bellow by  $\theta_1$ .
- 2. local similarity: There is a constant  $\mathcal{I}$  depending on  $\theta$  only such that for each  $1 \leq i \leq k$ ,  $el_{M_i} \leq \mathcal{I}el_{M_{i-1}}$ .

Therefore, the hierarchical gradient generated is  $(\theta_1, \mathcal{I}_1, |M_k|)$ -well-conditioned.

Theorem 3.8. Let  $M_1', ..., M_k'$  be any  $(\theta, \mathcal{I}, b)$ -well-conditioned gradient of  $M_0$ , for some positive constants  $\theta, \mathcal{I}$ . Let  $M_1, ..., M_k$  be the result of the application of the algorithm of Figure 3 on  $M_0$  for k iterations, then there is a constant c such that  $\forall i : |M_i| \leq c|M_i'|$ . Hence  $M_1, ..., M_k$  is  $(\theta_1, \mathcal{I}_1, c_1b)$ -well-conditioned gradient.

A simple corollary of the last theorem, is that our algorithm optimizes the number of levels up to an

**Procedure:** MULTI\_LEVEL\_COARSEN $(M_0)$ 

- 1. Let k be the length of the required hierarchy.
- 2. Let  $M_k = \text{ONE\_LEVEL\_COARSEN}(M_0, f_{C^{k+1}, P})$ . Let  $S_k$  be the point set of  $M_k$ .
- 3. For i = k 1 to 1
  - Let  $S_i$  be a maximal independent set of  $CG_i$  that contains  $S_{i+1}$ . Note that by Lemma 3.6  $S_{i+1}$  must be an independent set of  $CG_i$ .
  - Let  $M_i = DT(S_i)$ .
- 4. Return  $(M_1, ..., M_k)$ .

Figure 3: Multi level nested function based coarseing.

additive constant factor when coarseing the mesh down to a constant sized mesh:

Corollary 3.9. If  $M'_1,...,M'_k$  is  $(\theta,\mathcal{I},1)$ -well-conditioned gradient, then  $M_1,...,M_k$  can be completed to an  $(\theta_1,\mathcal{I}_1,1)$ -well-conditioned gradient of depth at most  $k+c_1$ .

# 4 Spacing Functions Qualities and Mesh Qualities

This section focuses on the intimate connection between the mesh qualities and its spacing function qualities. In particular, we show that spacing functions can capture the two most important mesh properties: its elements' shape, and its elements' size and number. This connection is the foundation of our coarsening approach correctness, which will be discussed in the next section.

**4.1 From a mesh to a spacing function.** The spacing function we recover from each mesh is the Nearest Neighbour function (NN), see Definition 3.2. This function depends only on the node set of the mesh. A generalization of this function, the local feature size function, plays an important role in the analysis of mesh generation algorithms [16, 3, 14]. These functions change slowly spatially, in analogy to the slowly changing element sizes of the well-shaped mesh.

Definition 4.1. A function f is 1-Lipschitz over a domain  $\Omega$  if for any two points x,y in  $\Omega$ ,  $|f(x)-f(y)| \leq ||x-y||$ .

The proofs of the following two lemmas are elementary: LEMMA 4.2. (RUPPERT [16]) For each point set P in  $\mathbb{R}^d$ ,  $NN_P$  is 1-Lipschitz.

Lemma 4.3.  $NN_P$  is a 2-spacing function of P.

For a good aspect ratio mesh M, the el function (see Definition 2.1) is equivalent up to a constant factor to NN.

Theorem 4.4. (Ruppert [16]) Let M be a mesh with smallest angle bound  $\theta$ . Let P be the node set of M. There exist two positive constants  $C_1, C_2$  depending on  $\theta$  only such that  $C_1 el_M(x) \leq NN_P(x) \leq C_2 el_M(x)$ .

Since every point set P is 2-spaced according to the 1-Lipschitz function  $NN_P$ , these notions by themselves are not powerfull enough to describe good aspect ratio meshes. Intuitively, spacing functions prevent points from clustering, but do not prevent the formation of arbitrarily large gaps. We now formalize the notion of a gap in terms of the spacing function.

Definition 4.5. Let f be an integrable function over a domain  $\Omega$ . The f-area of a sub-domain B of  $\Omega$  is given by  $A_f(B) = \int_B 1/(f(x))^2 dA$ .

The following lemma shows triangles of a good aspect ratio mesh are of constant *NN*-area, and relates the number of elements to the area of the domain. The proof is omitted from this extended abstract.

Lemma 4.6. Let M be a mesh with smallest angle bound  $\theta$ .

1. For each mesh triangle T:

$$\frac{\sin^{2}\theta}{2{C_{2}}^{2}} \leq A_{NN_{M}}\left(T\right) \leq \frac{1}{2C_{1}^{2}\sin\theta}$$

2. Let N be the number of triangles, then:

$$\frac{\sin^2 \theta}{2C_2^2} N \le A_{NN_M}(\Omega) \le \frac{1}{2C_1^2 \sin \theta} N$$

Where  $C_1$  and  $C_2$  are the constants of Theorem 4.4. We capture the notion of a gap using empty balls. A ball B is called an *empty* ball with respect to a mesh M if B does not contain any node of M. The following lemma states an empty ball can intersect at most a constant number of triangles. The proof is omitted from this paper.

Lemma 4.7. Let M be a mesh with smallest angle bound  $\theta$ . Any empty ball B intersects at most  $C_3 = 4 + 3 \frac{4}{\sin^4 \theta}$  triangles of M.

Finally, the following theorem states the relationship between a good aspect ratio mesh and spacing functions. In particular, a good aspect ratio mesh posseses a 2-spacing function and has no large gaps in terms of that spacing function. The next sub-section shows the inverse statement is true as well. Theorem 4.8 is a simple corollary of Lemmas 4.6 and 4.7.

THEOREM 4.8. Let M be a mesh with smallest angle bound  $\theta$ .

1. Let  $C_4 = C_3 \frac{1}{2 \sin \theta C_1^2}$ . The NN<sub>M</sub>-area of each empty ball is at most  $C_4$ .

2. Let  $L = 2 \cdot 8^{C_4}$ , and let x be any point of  $\Omega$ . Any ball containing x whose radius is greater than  $L \cdot NN_M(x)$  must contain some node p of the mesh.

The 1-dimensional analogs of the notion of gaps, and their related theorems, are valid for the function NN on the boundaries of  $\Omega$  (in our case the four edges of the unit square). They were omitted for the sake of brevity.

**4.2** From a spacing function to a mesh. In this section we show that a point set spaced according to a spacing function, where no large gaps in terms of the spacing function occur, is the node set of a good aspect ratio mesh. In particular, we show the Delaunay triangulation of that point set is of good aspect ratio.

The following theorem is the inverse of Theorem 4.8. Theorem 4.9. Let P be a set of points  $\beta$ -spaced by a 1-Lipschitz function f, such that each empty ball B is of constant bounded area,  $A_f(B) \leq \alpha$ . Then the smallest angle  $\theta$  of each Delaunay triangle of P is bounded by:  $\sin \theta \geq \frac{8^{-\alpha}}{2\beta}$ .

The following theorem shows that the function f  $\beta$ —spacing P is equivalent up to a constant factor to the function  $NN_P$ . The connection between f and  $el_{DT(p)}$  then follows by Theorem 4.4.

Theorem 4.10. Let P be a set of points  $\beta$ -spaced by a 1-Lipschitz function f, such that for each empty ball B  $A_f(B) \leq \alpha$ , then:

$$NN_P(x)/(L^2 + 2L) \le f(x) \le (2\beta + 1)NN_P(x)$$

For the sake of brevity, the theorems in this subsection ignored the boundary case. For the analogous results on the boundaries, we further require that the point set is spaced on the boundaries with no large gaps in terms of the spacing function restricted to the boundaries, and that points in the interior of  $\Omega$  stay away from the boundary: for some constant  $\delta$ , the distance of each non boundary point p from the boundary must be at least  $\delta \cdot f(p)$ . The aspect ratio of the Delaunay triangulation then depends on  $\delta$  as well.

### 5 Spacing Function Based Coarsening

This section is concerned with the correctness of our coarsening approach, as stated in Theorems 3.8 and 3.7. In particular, we show that given an initial bounded aspect ratio mesh and a coarsening factor C, our algorithm produces a bounded aspect ratio mesh which is (up to a constant factor) the smallest mesh whose el function is at most C times larger (point wise) than the el function of the original mesh.

**5.1** Aspect ratio of the coarser meshes. We first establish that our coarsening functions are all 1-

Lipschitz. This follows from Definitions 3.3 and 3.4.

LEMMA 5.1. If g is 1-Lipschitz, then for any coarsening factor  $C \geq 0$ ,  $f_{g,C,P}$  and  $t_{g,C,P}$  are 1-Lipschitz.

The next theorem establishes the main result of this sub-section: that the mesh generated by our one-level coarsening algorithm, see Figure 2, is a bounded aspect ratio mesh.

Theorem 5.2. Let P be the point set of a mesh  $M_0$  with smallest angle bound  $\theta$ . Let C>1 be the coarsening factor. Let  $M_1$  be the coarser mesh returned by our one-level coarsening algorithm with spacing parameter  $\beta=12\sqrt{L}$  and coarsening function  $f_{NN,C,P}$ . Further, assume  $L\geq 4$ , where L is the constant of Theorem 4.8. There exists a constant  $\theta_1$  depending on  $\beta$  and  $\theta$  only such that the smallest angle of  $M_1$  is larger than  $\theta_1$ . Proof: Let  $P_1$  denote the point set of  $M_1$ .  $P_1$  is  $\beta$ -spaced by  $f_{NN,C,P}$ . Theorem 4.9 implies that if the f-area of any Delaunay ball B is smaller than some positive constant  $\alpha$ , then  $\sin\theta_1\geq \frac{8^{-\alpha}}{2\beta}$ . Let f stand for  $f_{NN,\alpha,P}$ . Therefore, we proceed to prove the theorem by showing that the f-area of any empty ball in  $M_1$  is bounded .

We fix some point  $p \in P_1$  and a ball B through p. Without loss of generality, let f(p) = 1, let p = (0,0), and let the center of B lie on the y-axis, see Figure 4. We show that if the ball B is large, it cannot be empty. Let L be the constant of Theorem 4.8. Let B's radius be 12L. We show B must contain a point of  $P_1$  in its interior. (Note that B all ready contains p on its boundary).

The center coordinate of B is (0, 12L). Let  $B_0$  be a smaller ball nested in B of radius 3L centered at (0, 2+3L). The proof consists of showing that  $B_0$  must contain a point of P, and that this point could not be ruled out by taking a maximal independent set of the conflict graph of P, and hence must belong to  $P_1$  as well.

Since M is a good aspect ratio mesh, Theorem 4.8 implies any ball of radius  $LNN_P(x)$  through x must contain a point  $q \in P$ . Since  $f \geq NN_P$ , any ball of radius Lf(x) through x must contain a point  $q \in P$ . Let  $x_0 = (0, 2)$ .  $x_0$  lies on the boundary of  $B_0$ , and by the 1-Lipschitz property of f,  $f(x_0) \leq f(p) + 2 = 3$ . Hence, the ball  $B_0$  which is of radius  $3L \geq Lf(x_0)$  must contain a point  $q \in P$ .

Without loss of generality, let q = (x, y) be on the boundary of  $B_0$ . (all the bounds we derive hold if q is internal to  $B_0$  as well). We now show  $q \in P_1$  as well. We use the following facts about q:

I. By Lemma 5.3, the distance from q to the boundary of B is larger than y/2.

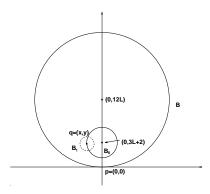


Figure 4: The existence of a mesh point of  $M_1$  in B.

II. By Lemma 5.4, if  $L \geq 4$  and  $\beta = 12\sqrt{L}$ , then  $f(q)/\beta \leq y/6$ .

Let  $B_1$  be the ball with center q and radius  $f/\beta$  (see the dashed ball in Figure 4.) If point  $q \notin P_1$ , then there must be some point  $w \in P_1$  such that q and w share an edge in the conflict graph used to create  $P_1$ . If  $w \in B$ then B contains a point of  $P_1$  in its interior, and we are done, therefore w must be outside B and property I implies  $d = ||w - q|| \ge y/2$ .

Because f is 1-Lipschitz,  $f(w) \leq d + f(q)$ . q and w share a conflict graph edge, hence  $f(q) + f(w) \geq \beta d$ . The last two inequalities imply  $2f(q) + d \geq \beta d$ . By property I,  $2f(q) \geq (\beta - 1)d \geq (\beta - 1)y/2$ . By property II,  $\beta y/3 \geq 2f(q) \geq (\beta - 1)y/2$ . Hence  $\beta/3 \geq (\beta - 1)/2$ . However, the last is a contradiction for  $\beta > 3$ , as is the case for our  $\beta$ .

This contradiction implies q and w do not share a conflict graph edge, and  $q \in P_1$ . This in turn implies that if B's radius is greater 12L, B can not be empty.

The following two lemmas were necessary for the proof of the theorem:

LEMMA 5.3. If L > 2/3 then the distance from any point (x, y) on  $B_0$  to the boundary of B is greater than y/2.

**Proof**: First notice that y>0. The distance from (x,y) to the boundary of B is equal to  $R-\sqrt{x^2+(y-R)^2}$ , where R=12L. Because (x,y) is on the boundary of  $B_0$ , we have  $x^2+(y-r-2)^2=r^2$ , implying  $x^2+y^2=2y(2+r)-4-4r=y(4+6L)-4-12L$ , for r=3L. This in turn implies, after some manipulations we omit, that  $x^2+(y-r-2)^2<(R-y/2)^2$ . This last inequality is true when L>2/3, y>0, L>0 and R=12L. Because y/2<R we have, after taking the square root,  $R-\sqrt{x^2+(y-R)^2}>y/2$ , completing the proof.

LEMMA 5.4. If L > 4 and  $\beta = 12\sqrt{L}$ , then

$$\frac{f_{NN,C,P}(q)}{\beta} \leq \frac{1 + \|p-q\|}{\beta} \leq \frac{y}{6}.$$

Proof:

As shown in the proof of Lemma 5.3  $\|p-q\|^2 = x^2 + y^2 = y(4+6L) - 4 - 12L$ . Thus,  $f(q) \le 1 + \|p-q\| \le 1 + \sqrt{y(4+6L) - 4 - 12L}$ . The Lemma states that  $1 + \sqrt{y(4+6L) - 4 - 12L} \le \beta y/6$ , or equivalently,  $\sqrt{y(4+6L) - 4 - 12L} \le \beta y/6 - 1$ 

Squaring both side, it suffices to show that  $y(4+6L)-4-12L<\beta^2y^2/36+1-\beta y/3$ . Hence it is enough to show that  $\beta^2y^2/36-\beta y/3-y(4+6L)>0$ . Since y>2 this is true if  $\beta^2/18-\beta/3-(4+6L)>0$ . The conditions L>4 and  $\beta=12\sqrt(L)$  guarantees this inequality.  $\square$ 

The aspect ratio bound derived in Theorem 5.2 applies to the special case of quasi-uniform meshes as well. However, we can obtain better bounds, using a simpler proof, for quasi-uniform meshes. These bounds are derived for a simpler coarsening spacing function, the threshold spacing function of Definition 3.4.

Theorem 5.5. Let P be the point set of a mesh  $M_0$  with smallest angle bound  $\theta$ . Let C>1 be the coarsening factor. Let  $M_1$  be the coarser mesh returned by our one-level coarsening algorithm with spacing parameter  $\beta=2$  and coarsening function  $t_{NN_P,C,P}$ . The smallest angle  $\theta_1$  of  $M_1$  is bounded bellow by:

$$\sin \theta_1 \ge \frac{1}{3 + 4L}$$

where L is the constant from Theorem 4.8, depending on  $\theta$  only.

**5.2** Local similarity. We now show that neighbouring meshes in the hierarchical coarsening gradient generated by our method our locally similar, see Definition 2.4.

The proof of the following lemma is elementary:

Lemma 5.6. [linearity and monotonicity] For any C > 1 and for any  $\gamma \geq 1$ . For each point  $x \in \Omega$ 

$$f_{g,C,P}(x) \le f_{g,\gamma C,P}(x) \le \gamma f_{g,C,P}(x).$$

Theorem 5.7. Let  $M_0$  be a mesh whose smallest angle bound is  $\theta$ . There exists a constant  $\mathcal{I}$  depending on  $\theta$  only such that such that for each mesh  $M_{i+1}$  of the hierarchical gradient  $(M_1, ..., M_k)$  produced by the Algorithm of Figure 3

$$el_{M_{i\perp 1}} \leq \mathcal{I}el_{M_i}$$

**Proof**: The coarsening function used to create  $M_{i+1}$  is  $f_{i+1} = f_{g,2C,P}$ , the one used to create  $M_i$  is  $f_i = f_{g,C,P}$ ,

for some value C. By Lemma 5.6,

$$f_{i+1} \leq 2f_i$$

By Theorem 4.10

$$NN_{P_i}/(L^2+2L) \le f_i \le (2\beta+1)NN_{P_i}$$

where  $P_j$  is the set of points of mesh  $M_j$ . By Theorem 4.4

$$C_1 el_{M_i} \leq NN_{P_i} \leq C_2 el_{M_i}$$

Hence

$$C_1/(L^2+2L)el_{M_{i+1}} \le f_{i+1} \le 2f_i$$

and

$$2f_i \le 2(2\beta + 1)C_2el_{M_i}$$

and we can take  $\mathcal{I}$  to be  $2C_2(2\beta+1)(L^2+2L)/C_1$ .  $\square$ 

**5.3 Size optimality.** We now show, up to a constant, that for any C > 1 the size of the mesh produced by our one-level coarsening algorithm is the smallest possible. The following Lemma shows that  $f_{g,C,P}$  is the largest 1-Lipschitz function that is smaller than  $C \cdot g$  at the points of the initial mesh  $M_0$ . The proof is elementary, and is omitted from this paper.

Lemma 5.8. Let C>1. Let h be a 1-Lipschitz function over the domain  $\Omega$  such that for all  $p\in P$   $h(p)\leq C\cdot g(p)$ . Then for all  $x\in\Omega$ ,  $h(x)\leq f_{g,C,P}(x)$ .

We now show that  $t_{NN_M,C,P}$  and  $f_{NN_M,C,P}$  are equivalent up to a constant factor for a quasi–uniform mesh M. This connection implies it suffices to show that meshes spaced by  $f_{NN_M,C,P}$  are of optimal size.

Lemma 5.9. Let M be a quasi-uniform mesh, i.e. there exist ratio constant  $\rho$  such that:

$$\min NN_M < NN_M(x) < \rho \min NN_M$$

Then

$$f_{NN_{M,C,P}}/(2L\rho + 2\rho) \le t_{NN_{M,C,P}} \le f_{NN_{M,C,P}}$$

Definition 5.10. Let M and M' be two well-shaped meshes over a domain  $\Omega$ . For any positive C>1, we say elements of M' are at most a factor of C larger than those of M if for all  $x\in\Omega$ ,  $el_{M'}(x)\leq C\cdot el_{M}(x)$ . In other words, for each point  $x\in\Omega$ , the largest triangle of M' that contains x is no more than a factor C larger than the largest triangle of M that contains x.

We now state the main result of this sub-section:

Theorem 5.11. Let  $M_0$  be mesh with smallest angle bound  $\theta_0$ . For any C>1, Let M be the mesh obtained by the algorithm of Figure 2 with coarsening factor C. Let M' be a mesh with minimal angle bound  $\theta'$  whose elements are at most a factor of C larger than those of  $M_0$ , then there exists a constant D depending on  $\theta_0$  and  $\theta'$  only such that  $|M| \leq D|M'|$ .

**Proof**: By the assumption  $el_{M'} \leq Cel_{M_0}$ . Since the constant bounds we derived in previous Lemmas depend on the smallest angle, we will refer to them in this proofs as a function of the relevant angle. Let  $P_0$  be the node set of  $M_0$  and let P be the node set of M. By Theorem 4.4,

$$\frac{NN_{M'}}{C_2(\theta')} \le el_{M'} \le Cel_{M_0} \le \frac{C}{C_1(\theta_0)} NN_{M_0}$$

Therefore by Lemma 5.8,

$$NN_{M'} \leq f_{NN_{M_0}, \frac{CC_2(\theta')}{C_1(\theta_0)}, P_0}$$

By Lemma 5.6:

$$NN_{M'} \le \frac{C_2(\theta')}{C_1(\theta_0)} f_{NN_{M_0,C,P_0}}$$

By Theorem 4.10:

$$NN_{M'} \le \frac{(2\beta(\theta_0) + 1)C_2(\theta')}{C_1(\theta_0)}NN_M$$

where  $\beta$  is the spacing constants used in the coarsening algorithm. Now Lemma 4.6 implies the result.

A simple application of this Lemma to all the levels of the coarsening gradient proves Theorem 3.8, whereas Theorem 5.2 can be applied to its multi-level version, Theorem 3.7.

# 6 Practical Concerns

**6.1 Mesh quality.** The proof outlined in previous sections provided a constant bound on the smallest angle of of the mesh hierarchy. However the bound can be quite small mathematically. Given the practical importance of the coarsening problem, we implemented the algorithm presented in this paper. We now provide some experimental evidence that our coarsening approach indeed produces coarsening meshes of very reasonable quality in practice.

We include the numerical data for the "crack plate" mesh, which was generated by Omar Ghattas and Xiaogang Li of Carnegie Mellon University. The physical problem modeled by the mesh is a plate with a horizontal crack running from the middle of the left edge to the center of the plate [8]. The following Table 1 lists the coarsening hierarchy statistics of our method. See Figure 5 for these meshes.

To offer a comparison, we also implemented a maximal independent set approach. We applied the MIS based program on the same initial mesh and observed a significant and iterative degradation on the smallest angle as shown in Table 2. This smallest angle occurs

coarsening	num	num	$\min$	max
factor	$\operatorname{nodes}$	triangles	angle	angle
1	5120	9984	41.25	93.88
2.1	2242	4374	18.43	135.01
4	1035	1982	18.23	135.06
8	305	564	18.43	135.01
16	148	272	19.44	123.69
32	119	220	18.01	130.23
64	102	189	18.01	130.23
128	92	170	18.01	130.23

Table 1: Quality of meshes coarsened using a function based approach. The first row describes the original mesh.

coarsening	num	num	min	max
iteration	$\operatorname{nodes}$	${ m triangles}$	angle	angle
1	1215	2346	18.43	126.87
2	289	537	14.04	139.40
3	76	131	3.37	135.00

Table 2: Quality of meshes coarsened using a maximal independent set approach.

in the center of the square, so the iterative degradation can not be attributed to edge effects, or our choice of boundary coarsening. The crack mesh is extremely unstructured, and the MIS based program fails to guarantee a bound on the aspect ratio in part because it attempts to reduce the number of mesh elements too much. Note that with our approach, the coarsening factor is doubled in later iterations but not the number of nodes.

**6.2** Algorithmic efficiency. The algorithm of Figure 2 has a simple  $O(n^2)$  time implementation. We an appropriate choice of data structures, we can reduce the algorithmic cost. The two computational expensive steps are the coarsened function construction and the point set selection according these functions.

The spacing function computation as described in Figure 2 takes  $O(n^2)$  time. In practice we can use other related spacing function, that still yield a theoretically correct algorithm. A particulary useful one is the spacing function based on a balanced quad-tree [3] of the fine mesh. The balanced quad-tree can be efficiently coarsened and can be used as an estimation of the spacing function of the underlying mesh. An important next step in our work is to develop such an efficient and practical implementation of our function based coarsening algorithm.

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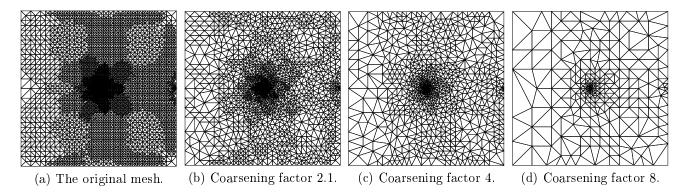


Figure 5: Coarsening the crack plate mesh using a coarsening function approach.

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