# 國立交通大學

# 應用數學系

數學建模與科學計算碩士班

# 碩士論文

Juli

多重網格與自調適法於 Laplace 方程角奇異解的 數值計算

Multigrid and Adaptive Methods for Computing Singular Solutions of Laplace Equation on Corner

Domains

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中華民國九十九年七月

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### 摘要

橢圓邊界值問題在凹角的地方會有奇異的行為,而這個奇異的行為對於 用有限元素法離散的精確度會受到影響。對於给定 Dirichlet 邊界條件的 Poisson 方程式和在定義域有凹角的情況之下,本論文利用一個奇異解的表 示法  $u=w+\sum_{j=1}^{J}\kappa_{j}s_{j}$ 算出較準確的近似值,其中  $\kappa_{j}=\frac{1}{\pi}\{\int_{\Omega}fs_{-1}dx+\int_{\Omega}u\Delta s_{-1}dx\}$ 在工 程上稱之為應力強度因子。這些量的精確計算在許多實際的工程問題上, 是一門很重要的課題。

# Multigrid and Adaptive Methods for Computing Singular Solutions of Laplace Equation on Corner Domains

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#### Abstract

Elliptic boundary value problems on domain with corners have singular behavior near the corners. Such singular behavior affect the accuracy of the finite element method throughout the whole domain. For the Poisson equation with homogeneous Dirichlet boundary conditions defined on a polygonal domain with re-entrant corners, it is well known that the solution has the singular function representation  $u=w+\sum_{j=1}^{j}\kappa_{j}s_{j}$ , where w is the regular part of the solution and  $s_{j}$  are known as singular functions that depend only on the corresponding re-entrant angles. Coefficients  $\kappa_{j}$  known as the stress intensity factors in the context of mechanics can be expressed in terms of u by extraction formula  $\kappa_{j} = \frac{1}{\pi} \{ \int_{\Omega} fs_{-1}dx + \int_{\Omega} u\Delta s_{-1}dx \}$ , where  $s_{-j}$  are known as dual singular function. Accurate calculation of these quantities is of great importance in many practical engineering problems. Similar singular function problems in [1, 2].

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# Contents

Introduction1							
1	Finite Ele	ement Method3					
	1.1 Intro	eduction of Finite Element Method					
	1.2 Vari	ational Formulation5					
	1.2.1	Existence and Uniqueness of Solution5					
	1.3 Finit	e Element Discretization					
	1.3.1	Linear Interpolation					
	1.3.2	Coordinate Transformations					
	1.3.3	Linear FEM Discretization					
	1.3.4	Partial Derivatives					
	1.3.5	Assembling the Element Matrix12					
	1.4 Erro	r Estimation12					
	1.4.1	Interpolation Error with Piecewise Linear Functions					
	1.4.2	A Priori Error Estimation					
	1.4.3	A Posterior Error Estimation and Adaptive Mesh-Refinement Techniques					
	1.5 Finit	e Element Approximation for Singular Functions21					
	1.5.1	Lagrange Interpolation					

	1.5.2 Singular Element	3				
2	Multigrid Method2	6				
	2.1 Introduction of Multigrid Method20	б				
	2.2 Relaxation Process	7				
	2.2.1 Jacobi Method	8				
	2.2.2 Gauss-Seidel Method	9				
	2.2.3 Successive Overrelaxation Method	0				
	2.3 Inter-grid Interpolation : Restriction and Prolongation	2				
	2.4 Multigrid Algorithm	4				
	2.5 Complexity					
	2.6 Numerical Experiments	7				
3	Research Method	3				
4	Numerical Results	9				
References						

# Introduction

The finite element method has become one of the most popular and effective methods for the numerical solution of partial differential equations, particularly for elliptic equations. In practice many important problems involve polygonal domains. Previously, the geometry of a problem would be restricted so that the triangulation elements fit the polygonal boundary exactly. Form a theoretical standpoint, under the assumption that solutions were sufficiently smooth, this case has been thoroughly analyzed. Unfortunately, in practice one is not likely to achieve the smoothness required for these previous analyses. It is the basic behavior of elliptic equations that solutions possess singularities at corners. These singularities substantially affect the rate of convergence of numerical approximations.

To handle this problem, here the two main procedures which have been proposed to overcome this difficulty. The first is based on mesh refinements and has been analysed by Babuska and kellogg [21],Raugel, Schatz and Wahlbin, Thatcher for instance. This method may be applied to most of the practical problems since it requires only a qualitative knowledge of the behaviour of the solution near the corners. The second consists in augmenting the space of trail funcitons in which one looks for the approximate solution. This is done by adding some of the singular solutions of the problem to the usual spaces of piecewise polynomial funcitons. For instance, S.C. Brenner and L.Y. SUNG [9], Babuska and Rosen-zweig [22], Kellogg, Lelievre, Djaoua and Ladeveze and Peyret.

In this thesis ,the first we introduce the singular element to capture the singular point at the corner and see the accuracy reduced. The advantage of the singular element is that there are many small tranguler near the singular point, the solution near the singular point can be approximation efficiency, but the disadvantage is that create large linear systems. To solve the large linear systems, we introduce the Multigrid method. The second, we applied the S.C. Brenner and L. Y. SUNG's method as a standard. The advantage of S.C. Brenner's method is that the stress intensity factors  $\kappa_j$  can be represent by the simple expression and correct calculation, but the disadvantage is that the lack of accuracy for the whole domain .We will improve the accuracy by introducing the adaptive mesh-refinement and adaptive cut-off function.

Finite element methods and their error estimation are given in section 1. Multigrid method are introduced in chapter 2. The Poisson equation and the singular function representation are given in chapter 3. Numerical results are carried out in chapter 4.

1896

# **Chapter 1 Finite Element Method**

## **1.1 Introduction of Finite Element Method**

The basic idea in any numerical method for solving a differential equation is first to discretize given continuous problem with infinite degrees of freedom to a discrete problem or with only finite degrees of freedom such that the differential equation is transformed into a system of linear equations which can be solved by using a computer.

Finite element method start from a reformulation of a given differential equation as an equivalent variational problem. In the case of elliptic equations this variational problem in basic case is a minimization problem of the form

Find 
$$u \in V$$
 such that  $F(u) \leq F(v)$  for all  $v \in V$  (1.1)

where V is a given set of admissible functions and  $F: V \to R$  is a functional. F(v) is the total energy associated with v and (1.1) corresponds to an equivalent characterization of the solution of the differential equation as the function in V that minimizes the total energy of the considered system. In general the dimension of V is infinite and thus in general the problem (1.1) can't be solved exactly. To obtain a problem that can be solved on a computer the idea in the finite element method is to replace V by a set  $V_h$  consisting of simple function only depending on finitely many parameters. This leads to a finitedimensional minimization problem of the form:

Find 
$$u_h \in V_h$$
 such that  $F(u_h) \leq F(v)$  for all  $v \in V_h$  (1.2)

This problem is equivalent to a linear or nonlinear system of equations. We hope that the solution  $u_h$  of this problem is sufficiently good approximation of the solution of the original minimization problem (1.1). Usually one chooses  $V_h$  to be a subset of V and in this case (1.2) corresponds to the classical Ritz-Galerkin method.

The advantage of finite element methods as compared with finite difference methods is that complicated geometry, general boundary conditions and variable or non-linear material properties can be handled relatively easily. In all these cases one meets unneccessary artificial complications with finite difference methodology. Further, the finite element method has a solid theoretical foundation which gives added reliability and in many cases makes it possible to mathematically analyze and estimate the error in the approximate finite element solution.

To solve a given differential or integral equation approximately using the finite element method, one has to go through basically the following steps:

- 1. variational formulation of the given problem
- 2. Mesh Generator
- 3. discretization using FEM: construction of the finite dimensional space  $V_h$  and choose basis function
- 4. Assemble the element matrix to obtain global matrix

From step  $1\sim4$ , one obtain a linear systems we will introduce at section 1.2.

## **1.2 Variational Formulation**

We will now consider the following boundary value problem for the Poisson equation:

$$\begin{cases} -\Delta u = f & \text{in } \Omega \\ u = g & \text{on } \partial \Omega \end{cases}$$
(1.3)

where  $\Omega$  is a bounded domain in the  $\mathbb{R}^2$  with boundary  $\partial \Omega, g$  is a constant, f is a given function, where

$$\Delta u = \frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2} \tag{1.4}$$

the equivalent variational problem is

$$-\int_{\Omega} (\Delta u) v dx = \int_{\Omega} f v dx \tag{1.5}$$

where v is test function in  $H^1_0(\Omega)$  ,  $v|_{\partial\Omega}=0$  .By taking integration by parts ,

1896

$$\int_{\Omega} \nabla v \nabla u dx = \int_{\Omega} v f dx + \overrightarrow{n} \cdot \nabla u v |_{\partial \Omega} = \int_{\Omega} v f dx \tag{1.6}$$

## 1.2.1 Existence and Uniqueness of Solution

Let a: $V \times V \rightarrow R$  be a bilinear mapping with following properties:

(1) a(.,.) is symmetric

(2) (Continuity) a(.,.) is continuous, i.e, there is a constant  $\beta > 0$  such that  $|a(v,w)| \le \beta ||v||_V ||w||_V \quad \forall v, w \in V$ 

(3) (Coercivity) a(.,.) is V-elliptic, i.e, there is a constant  $\alpha > 0$  such that  $a(v,v) \ge \alpha ||v||_V^2 \quad \forall v \in V$ 

(4) L is continuous, i.e., there is a constant  $\Lambda > 0$  such that  $|L(v)| \leq \Lambda ||v||_V \quad \forall v \in V$ 

**Theorem (Lax-Milgram theorem)** Let V be a Hilbert space with scalar product  $(., .)_V$ and corresponding norm  $|| \cdot ||_V$  (the V - norm). Suppose that a(., .) is a bilinear form on  $V \times V$  and L a linear form on V such that under the assumptions (1)-(4), there exists a unique  $u \in V$  such that

$$a(u,v) = L(v)$$
, for all  $v \in V$   
**1.3 Finite Element Discretization**

Let  $T_h = \{K\}$  is a triangulation of  $\Omega \subset R^2$ , the integral equation can be rewritten as

$$\sum_{K \in T_h(\Omega)} \int_K \nabla v \nabla u dx = \sum_{K \in T_h(\Omega)} \int_K v f dx$$
(1.7)

The finite element method is then employed to discretize the terms  $\int \nabla v \nabla u dx$  and  $\int v f dx$ on element, we first look the germetry on an element.

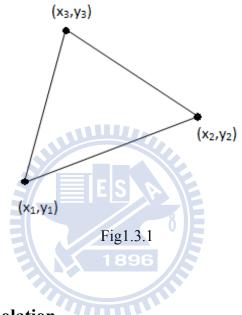
The geometry of the 3-node triangle is specified by the location of its three corner nodes on the  $\{x, y\}$  plane. The nodes are labeled 1, 2, 3 while traversing the sides in counterclockwise fashion. The location of the corners is defined by their coordinates:

$$x_i, y_i, i = 1, 2, 3$$

the area of triangle is denoted by A and is given by:

$$2A = \det \begin{bmatrix} 1 & 1 & 1 \\ x_1 & x_2 & x_3 \\ y_1 & y_2 & y_3 \end{bmatrix} = x_{21}y_{31} - x_{31}y_{21}$$

where  $x_{ij} = x_i - x_j$ ,  $y_{ij} = y_i - y_j$  for i, j = 1, 2, 3  $i \neq j$ .



#### **1.3.1** Linear Interpolation

One can choose a piecewise ploynomial function to approximate the exact solution u and the test function v. For example, if one choose linear piecewise function, then the function u(x, y) may be expressed as

$$u(x,y) = a_0 + a_1 x + a_2 y \tag{1.8}$$

where  $a_0$ ,  $a_1$  and  $a_2$  are coefficients to be determined from three conditions. In finite element work such conditions are often the nodal values taken by u at the corners:

 $u_1, u_2, u_3$ 

The expression in triangular coordinates makes direct use of these three values:

$$u(\zeta,\eta) = u_1(1-\zeta-\eta) + u_2\zeta + u_3\eta = \begin{bmatrix} u_1 & u_2 & u_3 \end{bmatrix} \begin{bmatrix} 1-\zeta-\eta \\ \zeta \\ \eta \end{bmatrix}$$
$$= \begin{bmatrix} 1-\zeta-\eta & \zeta & \eta \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \end{bmatrix}$$
(1.9)

equation (1.9) is called a linear interpolant for u.

## **1.3.2** Coordinate Transformations

Consider triangular on regular triangular, points of the triangle may also be located in terms

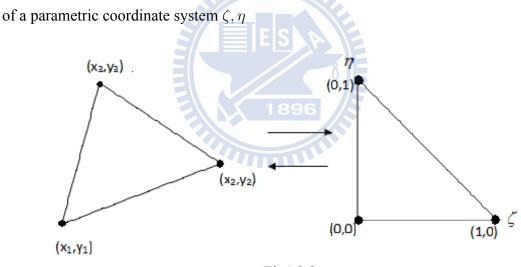


Fig1.3.2

Cartesian coordinates and triangular coordinates are linked by the relation

$$\begin{bmatrix} x \\ y \end{bmatrix} = \begin{bmatrix} x_1 \\ y_1 \end{bmatrix} [1 - \zeta - \eta] + \begin{bmatrix} x_2 \\ y_2 \end{bmatrix} \zeta + \begin{bmatrix} x_3 \\ y_3 \end{bmatrix} \eta$$
$$\begin{bmatrix} x_{21} & x_{31} \\ y_{21} & y_{31} \end{bmatrix} \begin{bmatrix} \zeta \\ \eta \end{bmatrix} + \begin{bmatrix} x_1 \\ y_1 \end{bmatrix}$$
(1.10)

These simply apply the linear interpolant formula to the Cartesian coordinates:  $x = x_1(1 - \zeta - \eta) + x_2\zeta + x_3\eta$  and  $y = y_1(1 - \zeta - \eta) + y_2\zeta + y_3\eta$ .

Inversion of (1.10) yields

$$\begin{bmatrix} \zeta \\ \eta \end{bmatrix} = \frac{1}{2A} \begin{bmatrix} y_{31} & x_{13} \\ y_{12} & x_{21} \end{bmatrix} \begin{bmatrix} x \\ y \end{bmatrix} - \frac{1}{2A} \begin{bmatrix} y_{31} & x_{13} \\ y_{12} & x_{21} \end{bmatrix} \begin{bmatrix} x_1 \\ y_1 \end{bmatrix}$$
(1.11)

#### **1.3.3 Linear FEM Discretization**

After generating mesh  $e_j$ , we shall now construct a finite dimensional subspace  $V_h$  of the space V defined above consisting of piecewise linear function. We now let  $V_h$  be the set of functions v such that v is linear on each subinterval  $e_j$ , v is continuous on domain  $\Omega$  and v = 0 on  $\partial\Omega$ . We observe that  $V_h \subset V$ . As parameter to describe a function  $u_j = v(x_j)$  we may choose the values  $u_j = v(x_j)$  at the node points  $x_j, j = 0, ..., m + 1$ . Let us introduction the basis function  $\phi_j \in V_h, j = 0, ..., m + 1$ .

$$\phi_j(x_i) = \begin{cases} 1 \text{ if } i = j \\ 0 \text{ if } i \neq j, \ i, j = 1, ..., M \end{cases}$$
(1.12)

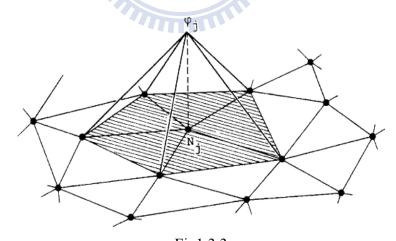


Fig1.3.3 i.e.,  $\phi_i$  is the continuous piecewise linear function that take the value 1 at node point

 $x_j$  and the value 0 at other node points. A function  $v \in V_h$  then has the representation

$$v(x) = \sum_{i=1}^{m} u_i \phi_i(x), \quad x \in \Omega$$
(1.13)

where  $u_j = v(x_j)$ , i.e., each  $u_j = v(x_j)$  can be written in a unique way as a linear combination of the basis function  $\phi_j$ . In particular it follow that  $V_h$  is a linear space of dimension m with basis  $\{\phi_j\}_{i=1}^m$ .

#### **1.3.4** Partial Derivatives

From equations (1.10) and (1.11) we immediately obtain the following relations between partial derivatives:

$$\frac{\partial x}{\partial \zeta} = x_{21}, \frac{\partial x}{\partial \eta} = x_{31}, \frac{\partial y}{\partial \zeta} = y_{21}, \frac{\partial y}{\partial \eta} = y_{31}.$$
$$\frac{\partial \zeta}{\partial x} = \frac{y_{31}}{2A}, \frac{\partial \eta}{\partial x} = \frac{y_{12}}{2A}, \frac{\partial \zeta}{\partial y} = \frac{x_{13}}{2A}, \frac{\partial \eta}{\partial y} = \frac{x_{21}}{2A}.$$

The derivatives of function  $u(\zeta, \eta)$  is

$$\nabla u = \begin{bmatrix} \frac{\partial u}{\partial x} \\ \frac{\partial u}{\partial y} \end{bmatrix} = \begin{bmatrix} \frac{\partial u}{\partial \zeta} \frac{\partial \zeta}{\partial x} + \frac{\partial u}{\partial \eta} \frac{\partial \eta}{\partial x} \\ \frac{\partial u}{\partial \zeta} \frac{\partial \zeta}{\partial y} + \frac{\partial u}{\partial \eta} \frac{\partial \eta}{\partial y} \end{bmatrix}$$

$$= \frac{1}{2A} \begin{bmatrix} \frac{\partial u}{\partial \zeta} y_{31} + \frac{\partial u}{\partial \eta} y_{12} \\ \frac{\partial u}{\partial \zeta} x_{13} + \frac{\partial u}{\partial \eta} x_{21} \end{bmatrix}$$

$$= \frac{1}{2A} \begin{bmatrix} y_{31} & y_{12} \\ x_{13} & x_{21} \end{bmatrix} \begin{bmatrix} \frac{\partial u}{\partial \zeta} \\ \frac{\partial u}{\partial \eta} \end{bmatrix}$$

$$= \frac{1}{2A} \Phi \begin{bmatrix} \frac{\partial u}{\partial \zeta} \\ \frac{\partial u}{\partial \eta} \end{bmatrix}$$
(1.14)

,where  $\Phi = \left[ egin{array}{cc} y_{31} & y_{12} \\ x_{13} & x_{21} \end{array} 
ight].$ 

Similarly, the derivatives of test function  $v(\zeta,\eta)$  is

$$\nabla v = \begin{bmatrix} \frac{\partial v}{\partial x} \\ \frac{\partial v}{\partial y} \end{bmatrix} = \frac{1}{2A} \begin{bmatrix} y_{31} & y_{12} \\ x_{13} & x_{21} \end{bmatrix} \begin{bmatrix} \frac{\partial v}{\partial \zeta} \\ \frac{\partial v}{\partial \eta} \end{bmatrix}$$
$$= \frac{1}{2A} \begin{bmatrix} \frac{\partial v}{\partial \zeta} & \frac{\partial v}{\partial \eta} \end{bmatrix} \Phi^{T}$$
(1.15)

So,we let

$$u = \sum_{i=1}^{3} u_i \phi_i \qquad , v = \sum_{i=1}^{3} v_i \phi_i$$
(1.16)

where  $\phi_1 = 1 - \zeta - \eta$ ,  $\phi_2 = \zeta$ ,  $\phi_3 = \eta$ ,

$$\begin{bmatrix} \frac{\partial u}{\partial \zeta} \\ \frac{\partial u}{\partial \eta} \end{bmatrix} = \begin{bmatrix} \frac{\partial \phi_1}{\partial \zeta} & \frac{\partial \phi_2}{\partial \zeta} & \frac{\partial \phi_3}{\partial \zeta} \\ \frac{\partial \phi_1}{\partial \eta} & \frac{\partial \phi_2}{\partial \eta} & \frac{\partial \phi_3}{\partial \eta} \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \end{bmatrix}$$
$$= \begin{bmatrix} -1 & 1 & 0 \\ -1 & 0 & 1 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \end{bmatrix} = D \begin{bmatrix} u_1 \\ u_2 \\ u_3 \end{bmatrix}$$
(1.17)

where  $D = \begin{bmatrix} -1 & 1 & 0 \\ -1 & 0 & 1 \end{bmatrix}$ . Similarity,  $\begin{bmatrix} \frac{\partial v}{\partial \zeta} \\ \frac{\partial v}{\partial \eta} \end{bmatrix} = \begin{bmatrix} v_1 & v_2 & v_3 \end{bmatrix} \begin{bmatrix} \frac{\partial \phi_1}{\partial \zeta} & \frac{\partial \phi_1}{\partial \eta} \\ \frac{\partial \phi_2}{\partial \zeta} & \frac{\partial \phi_2}{\partial \eta} \\ \frac{\partial \phi_3}{\partial \zeta} & \frac{\partial \phi_3}{\partial \eta} \end{bmatrix} = \begin{bmatrix} v_1 & v_2 & v_3 \end{bmatrix} D^T.$ for one element, we get

$$\int_{K} \nabla v \nabla u dx = \int_{\widehat{K}} \begin{bmatrix} v_1 & v_2 & v_3 \end{bmatrix} D^T \Phi^T \frac{1}{4A^2} \Phi D \begin{bmatrix} u_1 \\ u_2 \\ u_3 \end{bmatrix} |J| d\zeta d\eta \qquad (L.H.S)$$
(1.18)

 $K_K = \int_{\widehat{K}} D^T \Phi^T \frac{1}{4A^2} \Phi D |J| d\zeta d\eta$  is called element stiffness matrix.

and

$$\int_{K} v f dx = \int_{\widehat{K}} \begin{bmatrix} v_1 & v_2 & v_3 \end{bmatrix} \begin{bmatrix} \phi_1 \\ \phi_2 \\ \phi_3 \end{bmatrix} \begin{bmatrix} \phi_1 & \phi_2 & \phi_3 \end{bmatrix} \begin{bmatrix} f_1 \\ f_2 \\ f_3 \end{bmatrix} |J| d\zeta d\eta \ (R.H.S)$$
(1.19)

 $M_{K} = \int_{\widehat{K}} \begin{bmatrix} \phi_{1} \\ \phi_{2} \\ \phi_{3} \end{bmatrix} \begin{bmatrix} \phi_{1} & \phi_{2} & \phi_{3} \end{bmatrix} |J| d\zeta d\eta \text{ is called mass matrix.}$ 

#### **1.3.5** Assembling the element matrix

Assembling all the element matrices  $K_K$ ,  $M_k$  into global matrices, here  $\overrightarrow{u}, \overrightarrow{v}, \overrightarrow{f}$  are the global nodal vector. $\sum_{K \in T_h(\Omega)} K_K = K$ ,  $\sum_{K \in T_h(\Omega)} M_K = M$ . The equation (1.7) becomes

$$\overrightarrow{v} \cdot K \cdot \overrightarrow{u} = \overrightarrow{v} \cdot M \cdot \overrightarrow{f}$$
, for all  $v \in V_h$ 

As a result, we obtain a linear system

$$K \cdot \overrightarrow{u} = M \cdot \overrightarrow{f} \tag{1.20}$$

## **1.4 Error Estimation**

(Cea Lemma)

Let  $u \in V$  be the solution of a(u, v) = L(v),  $\forall v \in V$  and  $u_h \in V_h \subset V$ . Then

$$||u - u_h||_V \le \frac{\beta}{\alpha} \inf ||u - v_h||_V \quad \forall v_h \in V_h$$
(1.21)

where  $\alpha$  is coercivity scalar and  $\beta$  is continuity scalar.

For a typical elliptic problem satisfying the conditions (1)-(4) in the section 1.2.1,we choosing  $v_h = \pi_h u \in V_h$  to be a suitable interpolant of u and estimate the interpolation error  $||u - \pi_h u||_V$ .

#### **1.4.1** Interpolation Error with Piecewise Linear Functions

We first consider the case  $V = H^1(\Omega)$  and  $V_h = \{v_h \in V : v_h|_K \in P_1(K), \forall K \in T_h\}$ where  $T_h = \{K\}$  is a triangulation of  $\Omega \subset R^2$ , i.e.,  $V_h$  is the standard finite element space of piecewise linear functions on triangles K. For  $K \in T_h$  we define  $h_k$  = the dimeter of K=the longest side of K,

 $\rho_k =$  the dimeter of the circle inscribed in K,

$$h = \max_{K \in T_h} h_K.$$

We shall assume that there is a positive constant  $\beta$  independent of h, such that

$$\frac{\rho_K}{h_K} \ge \beta \qquad \forall K \in T_h \tag{1.22}$$

This condition means that the angles of the triangles K are not allowed to be arbitrarily small;the constant  $\beta$  is a measure of the smallest angle in any  $K \in T_h$ .Let  $N_i$ , i = 1, ..., M, be the nodes of  $T_h$ .Given  $u \in C^0(\Omega)$  we define the interpolant

$$\pi_h u(N_i) = u(N_i) \quad i = 1, ..., M$$
(1.23)

, thus  $\pi_h u$  is the piecewise linear function agreeing with u at the nodes of  $T_h$ .

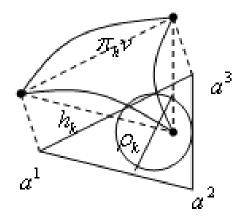


Fig1.4.1

For j=1,2 and  $x \in K$  we have

$$\sum_{i=1}^{3} \phi_i(x) = 1, \qquad (1.24)$$

$$\sum_{i=1}^{3} p_i(x)\phi_i(x) = 0, \qquad (1.25)$$

$$\sum_{i=1}^{3} \frac{\partial}{\partial x_j} \phi_i(x) = \frac{\partial}{\partial x_j} \sum_{i=1}^{3} \phi_i(x) = 0, \qquad (1.26)$$

$$\sum_{i=1}^{3} p_i(x) \frac{\partial \phi_i}{\partial x_j}(x) = \frac{\partial v}{\partial x_j}(x).$$
(1.27)

**Theorem 1** Let  $K \in T_h$  be a triangle with vertices  $a^i$ , i=1,2,3. Given  $v \in C^0(K)$ , let the interpolant  $\pi v \in P_1(K)$  be defined by

$$\pi v(a^i) = v(a^i), i = 1, 2, 3.$$
(1.28)

then

$$||v - \pi v||_{L^{\infty}(K)} \le 2h_{K}^{2} \max_{|\alpha=2|} ||D^{\alpha}v||_{L^{\infty}(K)},$$
(1.29)

$$\max_{|\alpha=1|} ||D^{\alpha}(v-\pi v)||_{L^{\infty}(K)} \le 6 \frac{h_{K}^{2}}{\rho_{K}} \max_{|\alpha=2|} ||D^{\alpha}v||_{L^{\infty}(K)},$$
(1.30)

where

$$||v||_{L^{\infty}(k)} = \max_{x \in K} |v(x)|.$$
(1.31)

**Proof.** Let  $\phi_i$ , i = 1, 2, 3, be the basis functions for  $P_1(K)$ . A general function  $w \in P_1(K)$ 

then has the representation

$$w(x) = \sum_{i=1}^{3} w(a^{i})\phi_{i}(x), \quad x \in K,$$
(1.32)

so that in particular

$$\pi v(x) = \sum_{i=1}^{3} v(a^{i})\phi_{i}(x), \quad x \in K,$$
(1.33)

since  $\pi v(a^i) = v(a^i)$ , we now using the Taylor expansion at  $x \in K$ :

$$v(y) = v(x) + \sum_{j=1}^{2} \frac{\partial v}{\partial x_j}(x)(y_j - x_j) + R(x, y),$$
(1.34)

where

$$R(x,y) = \frac{1}{2} \sum_{i,j=1}^{2} \frac{\partial^2 v}{\partial x_i \partial x_j} (\xi) (y_i - x_i) (y_j - x_j),$$
(1.35)

is the remainder term of order 2 and  $\xi$  is a point on line segment between x and y.In particular by choosing  $y = a^i$ , we have

$$v(a^{i}) = v(x) + p_{i}(x) + R_{i}(x)$$
(1.36)

where

$$p_{i}(x) = \sum_{j=1}^{2} \frac{\partial v}{\partial x_{j}}(x)(a_{j}^{i} - x_{j}), a^{i} = (a_{1}^{i}, a_{2}^{i}),$$

$$R_{i}(x) = R(x, a^{i}).$$
(1.37)

Since

$$|a_j^i - x_j| \le h_K, i = 1, 2, 3, \quad j = 1, 2,$$
(1.38)

the estimate of the remainder term  $R_i(x)$ 

$$R_i(x) \le 2h_K^2 \max_{|\alpha|=2} ||D^{\alpha}v||_{L^{\infty}(K)}, i = 1, 2, 3.$$
(1.39)

Now (1.33) and (1.36) combine to give

$$\pi v(x) = v(x) \sum_{i=1}^{3} \phi_i(x) + \sum_{i=1}^{3} p_i(x)\phi_i(x) + \sum_{i=1}^{3} R_i(x)\phi_i(x), \quad x \in K.$$
(1.40)

By(1.24), (1.25) in Lamma 2 and (1.40), we have

$$\pi v(x) = v(x) + \sum_{i=1}^{3} R_i(x)\phi_i(x), \qquad (1.41)$$

which gives us the following representation of the interpolation error:

$$v(x) - \pi v(x) = -\sum_{i=1}^{3} R_i(x)\phi_i(x).$$
(1.42)

Since  $0 \le \phi_i(x) \le 1$ , if  $x \in K$ , i = 1, 2, 3, we can use the previous estimate (1.39) of the

remainder term  $R_i$  to get

$$|v(x) - \pi v(x)| \leq \sum_{i=1}^{3} |R_i(x)| |\phi_i(x)|$$
  
$$\leq \max_i |R_i(x)| \sum_{i=1}^{3} |\phi_i(x)| \leq 2h_K^2 \max_{|\alpha|=2} ||D^{\alpha}v||_{L^{\infty}(K)}, x \in K(1.43)$$

we proves (1.29).

**Proof.** To prove (1.30) we differentiate (1.33) with respect to  $x_1$  to get

$$\frac{\partial \pi v}{\partial x_1}(x) = \sum_{i=1}^3 v(a^i) \frac{\partial \phi_i}{\partial x_1}(x), \qquad (1.44)$$

which together with (1.36) shows that

$$\frac{\partial \pi v}{\partial x_1}(x) = v(x) \sum_{i=1}^3 \frac{\partial \phi_i}{\partial x_1}(x) + \sum_{i=1}^3 p_i(x) \frac{\partial \phi_i}{\partial x_1}(x) + \sum_{i=1}^3 R_i(x) \frac{\partial \phi_i}{\partial x_1}(x).$$
(1.45)

15

Hence, by (1.26) and (1.27) we have

$$\frac{\partial \pi v}{\partial x_1}(x) = \frac{\partial v}{\partial x_1}(x) + \sum_{i=1}^3 R_i(x) \frac{\partial \phi_i}{\partial x_1}(x), \qquad (1.46)$$

which gives the following representation o the error  $\frac{\partial v}{\partial x_1} - \frac{\partial \pi v}{\partial x_1}$ :

$$\frac{\partial v}{\partial x_1}(x) - \frac{\partial \pi v}{\partial x_1}(x) = -\sum_{i=1}^3 R_i(x) \frac{\partial \phi_i}{\partial x_1}(x), \quad x \in K.$$
(1.47)

and

$$\max_{x \in K} \left| \frac{\partial \phi_i}{\partial x_1}(x) \right| \le \frac{1}{\rho_K},\tag{1.48}$$

which together with (1.39) finally gives

$$\left|\frac{\partial v}{\partial x_1}(x) - \frac{\partial \pi v}{\partial x_1}(x)\right| \le 6 \frac{h_K^2}{\rho_K} \max_{|\alpha=2|} ||D^{\alpha}v||_{L^{\infty}(K)}.$$
(1.49)

In the same way we estimate  $\frac{\partial v}{\partial x_2} - \frac{\partial \pi v}{\partial x_2}$  and thus (1.30) follows. The proof of the theorem is now complete once the lemma is established.

**Theorem 2** Under the assumptions of Theorem 1 there is an absolute constant C such that

$$||v - \pi v||_{L_2(K)} \le Ch_K^2 |v|_{H^2(K)}, \tag{1.50}$$

$$|v - \pi v|_{H^1(K)} \le C \frac{h_K^2}{\rho_K} |v|_{H^2(K)}.$$
(1.51)

$$||u - \pi_h u||_{H^1(\Omega)} \le Ch|u|_{H^2(\Omega)},$$
(1.52)

$$||u - \pi_h u||_{L^2(\Omega)} \le Ch^2 |u|_{H^2(\Omega)}$$
(1.53)

**Proof.** We have by summing over  $K \in T_h$ ,

$$\begin{aligned} ||u - \pi_h u||_{L^2(\Omega)}^2 &= \sum_{K \in T_h} ||u - \pi_h u||_{L^2(K)}^2 \leq \sum_{K \in T_h} C^2 h_K^4 |u|_{H^2(K)}^2 \\ &\leq C^2 h^4 \sum_{K \in T_h} |u|_{H^2(K)}^2 = C^2 h^4 |u|_{H^2(\Omega)}^2, \end{aligned}$$
(1.54)

and similarly using (1.22) , ie,  $\frac{h_K}{\rho_K} \leq \frac{1}{\beta},$ 

$$\begin{aligned} ||u - \pi_{h}u||_{H^{1}(\Omega)}^{2} &\leq \sum_{K \in T_{h}} C^{2} \frac{h_{K}^{4}}{\rho_{K}^{2}} |u|_{H^{2}(K)}^{2} \leq \sum_{K \in T_{h}} C^{2} \frac{h_{K}^{2}}{\beta^{2}} |u|_{H^{2}(K)}^{2} \\ &\leq C^{2} \frac{h^{2}}{\beta^{2}} |u|_{H^{2}(\Omega)}^{2} \end{aligned}$$
(1.55)

so that

$$||u - \pi_h u||_{H^1(\Omega)} \le \frac{Ch}{\beta} |u|_{H^2(\Omega)} = Ch|u|_{H^2(\Omega)},$$
(1.56)

if the constant  $\beta$  is included in the constant C,and

$$||u - \pi_h u||_{L^2(\Omega)} \le Ch^2 |u|_{H^2(\Omega)}.$$
(1.57)

# 1.4.2 A Priori Error Estimation

The bilinear form  $a(u, v) = \int \nabla u \nabla v dx$ , the finite element space of piecewise linear element is  $H^1(\Omega) = V$ , the test function space is  $H^1_0(\Omega)$ . The priori error estimate of Possion equation

$$||u - u_h||_{H^1(\Omega)} \le Ch|u|_{H^2(\Omega)}$$

Proof. by Céa Lamme, so the inequality is

$$||u - u_h||_{H^1(\Omega)} \le \frac{\beta}{\alpha} \inf ||u - v_h||_{H^1(\Omega)} \text{ for all } v_h \in V_h$$

where  $\alpha$  is coercivity scalar and  $\beta$  is continuity scalar ,than we choose  $v_h = \pi_h u$ , where  $\pi_h u$  is interpolant by the piecewise linear basis function

$$\frac{\beta}{\alpha}\inf||u-v_h||_{H^1(\Omega)} \le \frac{\beta}{\alpha}||u-\pi_h u||_{H^1(\Omega)}$$

by interpolation error estimate

$$\frac{\beta}{\alpha}||u-\pi_h u||_{H^1(\Omega)} \le C\frac{\beta}{\alpha}h|u|_{H^2(\Omega)} = C_2h|u|_{H^2(\Omega)}$$

## **1.4.3 A Posterior Error Estimation and Adaptive Mesh-Refinement** Techniques

On the other way to refine the mesh ,we introduce a posterior error estimation and the adaptive mesh-refinement techniques. Before presenting the a posteriori error estimators, we introduce some notations.

For any open subset  $\omega$  of  $\Omega$  with Lipschitz boundary  $\gamma$ , we denote by  $L^2(\omega)$ ,  $H^k(\omega)$ , and  $L^2(\gamma)$ ,  $k \ge 1$ , the standard Lebesque and Sobolev spaces ,respectively, equipped with the norms  $|| \cdot ||_{0,\omega} := || \cdot ||_{L^2(\omega)}, || \cdot ||_{k,\omega} := || \cdot ||_{H^k(\omega)}, || \cdot ||_{0,\gamma} := || \cdot ||_{L^2(\gamma)}.$ 

For  $T \in \overline{T}_h$  we denote by  $\overline{E}(T)$  and  $\overline{N}(T)$  the set of its edges and vertices, respectively. Let

$$\overline{E}_h := \bigcup_{T \in \overline{T}_h} \overline{E}(T), \qquad \overline{N}_h := \bigcup_{T \in \overline{T}_h} \overline{N}(T),$$

be the set of all edges and vertices, respectively, in the triangulation. We split  $\overline{E}_h$  and  $\overline{N}_h$  in the form

$$\overline{E}_{h} = \overline{E}_{h,\Omega} \bigcup \overline{E}_{h,D}, \qquad \overline{N}_{h} = \overline{N}_{h,\Omega} \bigcup \overline{N}_{h,D},$$

with

$$E_{h,D} := \{E \in E_h : E \subset \Gamma_D\}$$
  
$$\overline{N}_{h,D} := \{x \in \overline{N}_h : x \subset \Gamma_D\}$$

For  $T \in \overline{T}_h$  and  $E \in \overline{E}_h$ , we denote by  $h_T$  and  $h_E$  their diamter and length, respectively. For  $T \in \overline{T}_h$  and  $E \in \overline{E}_h$  and  $x \in \overline{N}_h$ , let

$$\omega_T := \bigcup_{\overline{E}(T) \cap \overline{E}(T') \neq 0} T', \ \omega_E := \bigcup_{E \in \overline{E}(T')} T', \ \omega_x := \bigcup_{x \in \overline{N}(T')} (T'),$$

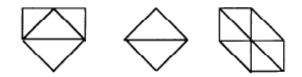


Figure 1.4.2: Domain  $\omega_T, \omega_E$  and  $\omega_x$ 

and put

$$V_T := \{\phi \in C(T) : \phi \in \prod_{\max(k+1,3),2}, \phi|_E \in \prod_{k+1,1}, \forall E \in \overline{E}(T) \quad (1.58)$$
  
$$\phi(x) = 0, \forall x \in \overline{N}(T), \phi|_E = 0, \forall E \in \overline{E}(T) \cap \overline{E}_{h,D}\},$$

$$V_x := \{ \phi \in C(\overline{\omega}_x) : \phi|_{T'} \in V_{T'}, \forall T' \subset \omega_x \}.$$
(1.59)

Given  $E \in \overline{E}_{h,\Omega}$  and  $\phi \in L^2(\omega_E)$  with  $\phi|_{T'} \in C(T'), \forall T' \subset \omega_E$ , we denote by  $[\phi]_E$  the jump of  $\phi$  across E in an arbitrary, but fixed direction. Put

$$R_E(u_h) := \begin{cases} -\left[\frac{\partial u_h}{\partial n}\right]_E, & \forall E \in \overline{E}_{h,\Omega}, \\ 0, 1896 & \forall E \in \overline{E}_{h,D}. \end{cases}$$
(1.60)

and

$$R_T := \prod_T f + \Delta u_h, \ \forall T \in \overline{T}_h,$$
(1.61)

where  $\prod_T$  is the  $L^2$  projectors of  $L^2(\Omega)$  onto the space of piecewise constant functions with respect to  $\overline{T}_h$ .

The first error estimator simply is a weighted combination of the residuals  $R_T(u_h)$ and  $R_E(u_h)$ . It is given by

$$\eta_{T,R} := \{h_T^2 || R_T(u_h) ||_{0,T}^2 + \sum_{E \in \overline{E}(T) \setminus \overline{E}_{h,D}} h_E || R_E(u_h) ||_{0,E}^2 \}^{1/2}, \ \forall T \in \overline{T}_h$$
(1.62)

The second error estimator is based on the solution of local Dirichlet problems. For any  $x \in \overline{N}_{h,\Omega}$ , it is given by

$$\eta_{x,D} := ||\nabla (u_{x,D} - u_h)||_{0,\omega} , \qquad (1.63)$$

For more detialed proof of the lower bound and upper bound for the error estimators is in [17].

According to the error estimator  $\eta_T$  we obtained above, we want to decide a refinement of T, there is a strategy that we refine all T with  $\eta_T \ge \gamma \max_{T \in \overline{T}_h} \eta_T$ . Here  $0 \le \gamma \le 1$ is a given threshold,typically  $\gamma = 0.5$ . This strategy is very cheap and often satisfactory results.

Having decided which elements should be refined, we refine them by connecting the mid-points of their edges. Triangles may be cut into four new ones by connecting the mid-points of their edges .

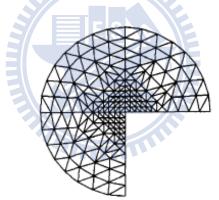


Figure 1.4.3: triangles with adaptive mesh-refinement

## **1.5** Finite Element Approximation for Singular Funcitons

The problem on hand is to devise singular elements within which the field variable would vary as  $R^t (0 \le t \le 1)$  where R is the radial distance from the point where the derivative of the function would have a singularity of the type  $R^{t-1}$ . For example, let  $(x_1, y_1)$  be the desired points of singularity in two-dimensional problem, such that R is equal to  $[(x-x_1)^2 +$   $(y - y_1)^2$ <sup>1/2</sup>. The singularity may be introduced in the geometric transformation between (x, y) and  $(\zeta, \eta)$  systems.

#### **1.5.1 Lagrange Interpolation**

For one-dimensional domain, consider a straight line along x-axis in a Cartesain reference, such that  $x_0 \le x \le x_M$ .Consider the dimensionless coordinate  $\xi = (x - x_0)/(x_M - x_0)$ such that  $0 \le \xi \le 1$ .A mth-order polynomial  $F^{(m)}(\xi)$  can be represented by interpolating the discrete values of the function at (m + 1) equidistant points  $\xi_i[i = 0, ..., m; \xi_i(i/m)]$ using the Lagrange interpolation function,  $L_i^{(m)}(\xi)$ , as

$$F^{(m)}(\xi) = \sum_{i=0}^{m} F_i L_i^{(m)}(\xi)$$
(1.64)

where

$$L_{i}^{(m)}(\xi) = \prod_{j=0, j \neq i}^{m} \left(\frac{\xi - \xi_{j}}{\xi_{i} - \xi_{j}}\right).$$
(1.65)

It is seen that the functions in (1.65) have the property that

$$L_i^{(m)}(\xi_k) = \delta_{ik} \tag{1.66}$$

where  $\delta_{ik} = 0$   $(i \neq k)$  and  $\delta_{ik} = 1(i = k)$ .

For a simple example, let we consider second-order polynomial  $F^{(2)}(\xi)$  and the transformation reduces to

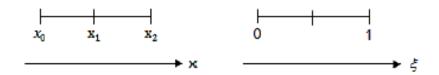


Fig1.5.1: x and  $\xi$  coordinate systems

$$(x - x_0) = (x_M - x_0)\xi \tag{1.67}$$

the Lagrange function can be represent as

$$L_0^{(2)} = 2(\xi - \frac{1}{2})(\xi - 1)$$
$$L_1^{(2)} = -4\xi(\xi - 1)$$
$$L_2^{(2)} = 2\xi(\xi - \frac{1}{2})$$

# 1.5.2 Singular Element

Apart from problems of fracture, at a given point in the domain, say P, the solution funciton F, whould have a singularity in its derivative. If R is the radial length measured from the point P, the solution function F would behave as  $R^{\lambda}$  ( $0 < \lambda < 1$ ), so that the first derivative of F at P would be infinite. Such singularities may arise at re-entrant corners, tips of sharp crack, etc.

For One-dimensional cases, consider a straight line along the x-axis,  $x_0 \le x \le x_M$ . Let the geometry of this line be described by an rth-order polynomial as

$$x = \sum_{i=0}^{r} x_i L_i^{(r)}(\xi)$$
(1.68)

where  $x_i = x(\xi_i)$  and  $\xi_i = (i/r)(i = 0, 1, ..., r)$  are (r + 1) equidistant points in the parametric domain  $0 \le \xi \le 1.(x_i$  themselves may not necessarily be equidistant). Also let the dependent variable, say u, be represented by an mth-order polynomial.

$$u(\xi) = \sum_{j=0}^{m} u_j L_j^{(m)}(\xi)$$
(1.69)

where  $u_j = u(\xi_j)$ , but now  $\xi_j = (j/m)(j = 0, ..., m)$  are (m + 1) equidistant points in  $0 \le \xi \le 1$ .

If the geometric transformation is defined by

$$(x - x_0) = (x_M - x_0)\xi^t \tag{1.70}$$

where t is a positive integer greater than unity, the function u(x) corresponding to  $u(\xi)$ of Eq.(1.69) would be an mth-order polynomial in the variable  $(x - x_0)^{1/t}$  or  $R^{1/t}$ , and hence the derivative  $(\partial u/\partial x)$  has a term which varies as  $R^{(1-t)/t}$ , for t > 1, would have a singularity at  $x = x_0$ .

For instance when r = 2 and t = 2, the positioning of nodes at  $x = x_0; x = x_0 + (x_M - x_0)(\frac{1}{2})^2 = x_0 + (x_M - x_0)(\frac{1}{4})$ , and  $x = x_M$  it would lead to  $x \sim \xi^2$  and hence  $\partial u / \partial x$  would have a singularity of the type  $R^{-1/2}$ .

In general, singularities in  $(\partial u/\partial x)$  of the type  $R^{(1-t)/t}$ ,  $t \leq r$  (t and r are integer), can be created by suitably choosing the valus  $x_i$  in an rth-order geometric transformation of the type (1.68).

For two-dimensional domain, consider a standard dimensionless triangle  $(0 \le \eta, \zeta \le 1)$  as shown in Fig1.5.2. The triangle  $P_1P_2P_3$  with Cartesain coordinate  $(x_i, y_i)(i = 1, 2, 3)$  is then mapped into the standard triangle ,through the relations:

$$(x - x_1) = (x_2 - x_1)\zeta + (x_3 - x_1)\eta$$
  

$$(y - y_1) = (y_2 - y_1)\zeta + (y_3 - y_1)\eta$$
(1.71)

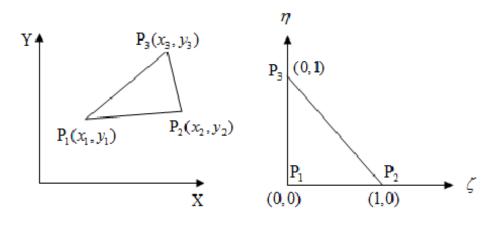


Fig1.5.2: Triangular element in (x, y) and  $(\zeta, \eta)$  coordinate systems Consider (m + 1) equidistant point along the line  $\eta = 0$  and  $\zeta = 0$ , respectively, such

that  $\eta_i = (i/m)$  and  $\zeta_j = (j/m)$  (i, j = 0, 1, ..., m). A mth-order polynomial (m being an integer),  $F^{(m)}(\eta, \zeta)$ , can be represented by interpolating the values of the function in the standard triangle, as:

$$F^{(m)}(\eta,\zeta) = \sum_{i=0}^{m} \sum_{j=0}^{m-i} F_{ij} L_{ij}^{(m)}(\eta,\zeta)$$
(1.72)

where

$$L_{ij}^{(m)} = \left[\prod_{r=0}^{i-1} \left(\frac{\eta - \eta_r}{\eta_i - \eta_r}\right)\right] \left[\prod_{s=0}^{j-1} \left(\frac{\zeta - \zeta_s}{\zeta_j - \zeta_s}\right)\right] \left[\prod_{t=0}^{m-(i+j+1)} \left(\frac{1 - (\eta + \zeta) - (t/m)}{1 - (\eta_i + \zeta_j) - (t/m)}\right)\right].$$
(1.73)

The function also have the property that

$$L_{ij}^{(m)}(\eta_k,\zeta_l) = \delta_{ik}\delta_{jl}.$$
(1.74)

For the detail of singular element geometric transformation in two-dimensional domain, see the reference [8].

# **Chapter 2 Multigrid Method**

## 2.1 Introduction of Multigrid Method

Multigrid methods(MG) are among the most efficient methods of solving the linear systems arising from discretization of eilliptic partial differential equations. There has been intensive research on the convergence of MG since it was introduced by Fedorenko. For symmetric positive-definite eilliptic problems, thanks to many researchers, such as Bank, Brandt, Braess, Bramble, Dupont, Hackbusch, Mandel and McCormick, etc, the convergence theory has matured. The major ingredients for convergence analysis of MG are called the approximation property and the smoothing property. One approach for convergence analysis is the so-called *compact perturbation technique*, which relies on a strong approximation property and treats the lower order terms as a small perturbation of the symmetric positive define term. The technique has been successfully applied to diffusion problems and Bramble, Pasciak, Wang, Xu have shown robust MG uniform convergence. In these studies, uniform convergence of MG can be established with one step of standard Jacobi or Gauss-Seidel smoothing even without regularity assumptions. Another approach requires a strong smoothing property to compensate for poor approximation property. In this direction, it is very important to find a robust smoother. Robust smoothers such as the block Jacobi, block Gauss-Seidel method and the incomplete LU factorization (ILU) method are commonly used.

The efficiency of the multigrid algorithm is achieved from an elegant combination of the smoothing procedure and the coarse grid correction procedure. The smoothing procedure plays the role of reducing highly oscillatory error modes, and the coarse grid is used to correct the remaining smooth error modes. Hackbusch and Braess give the first rigorous proof on the multigrid convergence and identify that the *smoothing property* and the *approximation property* are the cornerstones for the convergence analysis of multigrid methods. The smoothing property achieved by stationary iterative method. The approximation property is achieved by choosed intergrid interpolation operator.

In general, the stationary iterative methods can effective eliminating the high-frequency or oscillatory components of the error, but leave the low-frequency or smooth components of the error as we will shown in experiment 1 in section 2.6. This process is called the relaxation process in MG. Moreover, the smooth error modes on a fine grid appear more oscillatory on the coarse grid. As a result, repeating the relaxation process may consequently remove various oscillartory modes of error. To achieve the efficiency of the MG algorithm, we only have to concern how to restrict smooth error to the coarse grid and how to prolong the errors remained in coarse grid to the fine grid. This process is called the inter-grid interpolation, we will introduce the relaxation and interpolation process in the fillowing sections.

## 2.2 Relaxation Process

For relaxation process, we introduce the stationary iteration methods. Stationary iterative methods solve a linear system with an operator approximating the original one and based

on a measurement of the error in the result (the residual), form a "correction equation" for which this process is repeated. While these methods are simple to derive, implement, and analyze, convergence is only guaranteed for a limited class of matrices. Examples of stationary iterative methods are the Jacobi method, Gauss–Seidel method and the Successive over-relaxation method.

#### 2.2.1 Jacobi Mehtod

One of the methods to solve the linear system is the Jacobi method . The Jacobi method is derived by examining each of the N-1 equations in the linear system Av = f in isolation . If in the i - th equation

$$\sum_{j=1}^{N-1} a_{i,j} v_j = f_i,$$
(2.75)

we solve for the value of  $v_i$  while assuming the other entries of v remain fixed ,we obtain

$$v_i = (f_i - \sum_{j \neq i} a_{i,j} v_j) / a_{i,i} \quad , 1 \le j \le N - 1$$
(2.76)

this suggests an iterative method defined by

$$v_i^{(k)} = (f_i - \sum_{j \neq i} a_{i,j} v_j^{(k-1)}) / a_{i,i} \quad , 1 \le j \le N - 1$$
(2.77)

which is the Jacobi method.

In matrix terms, the definition of the Jacobi method can be expressed as

$$v^{(k)} = D^{-1}(L+U)v^{(k-1)} + D^{-1}f,$$
(2.78)

where the matrices D, -L and -U represent the diagonal, the strictly lower-triangular, and the strictly upper-triangular parts of A, respectively. The pseudocode for the Jacobi method is given by following

```
Choose an initial guess v^{(0)} to the solution u
for k = 1, 2, \cdots
for i = 1, 2, \cdots, N - 1
v_i = 0
for j = 1, 2, \cdots, i - 1, i + 1, \cdots, N - 1
v_i = v_i + a_{i,j}v_j^{(k-1)}
end
v_i = (f_i - v_i)/a_{i,i}
end
v^{(k)} = v
check convergence; continue if necessary
end
```

There is a simple modification which can be made to the Jacobi iteration.

$$v_j^* = (f_i - \sum_{j \neq i} a_{i,j} v_j^{(k-1)}) / a_{i,i} \quad , 1 \le j \le N - 1$$

However,  $v_j^*$  is now only an intermediate value. The new approximation is given by the weighted average

$$v_j^{(k)} = (1 - \omega)v_j^{(k-1)} + \omega v_j^* , 1 \le j \le N - 1$$

where  $\omega \in \mathbb{R}$  is a weighting factor which may be chosen. This iteration is called weighted Jacobi method.

#### 2.2.2 Gauss-Seidel Method

Consider again the linear systems Av = f, now assume that the equations are examined one at a time in sequence, and that previously computed results are used as soon as they are available ,we obtain the Gauss-Seidel method:

$$v_i^{(k)} = (f_i - \sum_{j < i} a_{i,j} v_j^{(k)} - \sum_{j > i} a_{i,j} v_j^{(k-1)}) / a_{i,i} \quad , 1 \le j \le N - 1$$
(2.79)

In matrix form,

$$Dv^{(k)} - Lv^{(k)} - Uv^{(k-1)} = f$$

$$(D-L)v^{(k)} = Uv^{(k-1)} + f$$

$$v^{(k)} = (D - L)^{-1}Uv^{(k-1)} + (D - L)^{-1}f$$

where the matrices D, -L and -U represent the diagonal, the strictly lower-triangular, and the strictly upper-triangular parts of A, respectively. The pseudocode for the Gauss-Seidel method is given by following

Choose an initial guess 
$$v^{(0)}$$
 to the solution  $u$   
for  $k = 1, 2, \cdots$   
for  $i = 1, 2, \cdots, N - 1$   
 $v_i = 0$   
for  $j = 1, 2, \cdots, i - 1$   
 $v_i = v_i + a_{i,j}v_j^{(k)}$   
end  
for  $j = i + 1, \cdots, N - 1$   
 $v_i = v_i + a_{i,j}v_j^{(k-1)}$   
end  
 $v_i = (f_i - v_i)/a_{i,i}$   
end  
 $v^{(k)} = v$   
check convergence; continue if necessary  
end

### 2.2.3 Successive Overrelaxation Method

The Successive Overrelaxation Method, or SOR ,is devised by applying extrapolation to the Gauss-Seidel method. This extrapolation takes the form of a weighted average between the

previous iterate and the computed Gauss-Seidel iterate successively for each component:

$$v_i^{(k)} = \omega \overline{v_i}^{(k)} + (1 - \omega) v_i^{(k-1)}$$
(2.80)

,where  $\overline{v_i}^{(k)}$  denotes a Gauss-Seidel iterate ,and  $\omega$  is the extrapolation (weighting) factor. The idea is to choose a value for  $\omega$  that will accelerate the rate of convergence of the iterates to the solution.

In matrix form, the SOR is written as

$$Dv^{(k)} = \omega D\overline{v}^{(k)} + (1 - \omega)Dv^{(k-1)}$$
  
=  $\omega [f + Lv^{(k)} + Uv^{(k-1)}] + (1 - \omega)Dv^{(k-1)}$ 

$$\implies (D - \omega L)v^{(k)} = [\omega U + (1 - \omega)D]v^{(k-1)} + \omega f$$
$$v^{(k)} = (D - \omega L)^{-1}[\omega U + (1 - \omega)D]v^{(k-1)} + (D - \omega L)^{-1}\omega f$$

where the matrices D, -L and -U represent the diagonal, the strictly lower-triangular, and the strictly upper-triangular parts of A, respectively.

If the extrapolation factor  $\omega$  is choosing by one,the SOR method simplifies to the Gauss-Seidel method. If  $0 < \omega < 1$ , its called underrelaxation. If  $1 < \omega < 2$ , its called overrelaxation and SOR fails to converge when  $\omega$  is outside the interval (0, 2).

The pseudocode for the SOR method is given by following

```
Choose an initial guess v^{(0)} to the solution u
for k = 1, 2, \cdots
for i = 1, 2, \cdots, N - 1
\sigma = 0
for j = 1, 2, \cdots, i - 1
\sigma = \sigma + a_{i,j}v_j^{(k)}
end
for j = i + 1, \cdots, N - 1
\sigma = \sigma + a_{i,j}v_j^{(k-1)}
end
\sigma = (f_i - \sigma)/a_{i,i}
v_i^{(k)} = v_i^{(k-1)} + \omega(\sigma - v_i^{(k-1)})
end
check convergence; continue if necessary
end
```

## 2.3 Inter-grid Interpolation : Restriction and Prolongation

First, we consider the linear prolongation, the operator will be denoted  $I_{2h}^h$ . It takes coarse grid vectors and produces find grid vectors according to the rule  $I_{2h}^h v^{2h} = v^h$ 

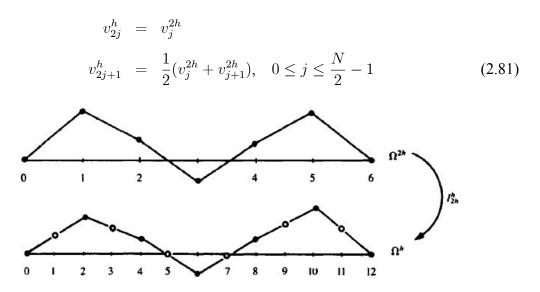


Figure 2.3.1: Prolongation of a vector on the coarse grid to the fine grid.

At even-numbered fine grid points ,the values of the vector are transferred directly from  $\Omega^{2h}$  to  $\Omega^h$ . At odd-numbered find grid points,the value of the vector is the average of the adjacent coarse grid values.

For the case N = 8, this operator has the form

$$I_{2h}^{h}v^{2h} = \frac{1}{2} \begin{bmatrix} 1 & & \\ 2 & & \\ 1 & 1 & & \\ & 2 & \\ & 1 & 1 & \\ & & 2 & \\ & & 1 & \end{bmatrix} \begin{bmatrix} v_1 \\ v_2 \\ v_3 \end{bmatrix}_{2h} = \begin{bmatrix} v_1 \\ v_2 \\ v_3 \\ v_4 \\ v_5 \\ v_6 \\ v_7 \end{bmatrix}_h = v^h$$

The second ,we consider moving vectors from a fine grid to a coarse grid. There are called restriction operators and are denoted by  $I_h^{2h}$ . It is defined by  $I_h^{2h}v^h = v^{2h}$ , where

$$v_j^{2h} = \frac{1}{4}(v_{2j-1}^h + 2v_{2j}^h + v_{2j+1}^h), \quad 1 \le j \le \frac{N}{2} - 1.$$
 (2.82)

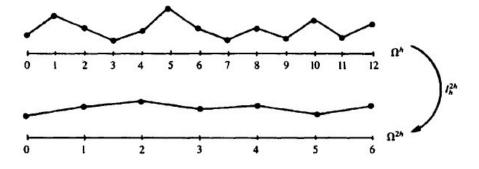


Figure 2.3.2: Restriction by full weighting of a find grid vector to the coarse grid.

the values of the coarse grid vector are a weighted average of values at neighboring find grid points.

In the case of N = 8, the operator has the form

$$I_{h}^{2h}v^{h} = \frac{1}{4} \begin{bmatrix} 1 & 2 & 1 & & \\ & 1 & 2 & 1 & \\ & & & 1 & 2 & 1 \\ & & & & 1 & 2 & 1 \end{bmatrix} \begin{bmatrix} v_{1} \\ v_{2} \\ v_{3} \\ v_{4} \\ v_{5} \\ v_{6} \\ v_{7} \end{bmatrix}_{h} = \begin{bmatrix} v_{1} \\ v_{2} \\ v_{3} \end{bmatrix}_{2h} = v^{2h}.$$

## 2.4 Multigrid Algorithm

After kowning the procedure, we compose all the components above to make the correction scheme. First, relaxation on the fine grid will eliminate the oscillatory components of the error, leaving a relatively smooth error, then transfer the residual on the fine grid to the coarse grid an solve the residual equation exactly on the  $\Omega^{2h}$ . Since the error is smooth on the  $\Omega^{2h}$ , we can prolongation the error accurately back to the fine grid. Finally, relaxation on the fine grid use the better initial which is correction before.

The procedure is given by following

 $\begin{array}{l} v^h \leftarrow MG(v^h, f^h) \\ \text{Relax } \nu_1 \text{ times on } A^h u^h = f^h \text{ on } \Omega^h \text{ with initial guess } v^h. \\ \text{Compute } r^{2h} = I_h^{2h}(f^h - A^h v^h). \\ \text{Solve } A^{2h}e^{2h} = r^{2h} \text{ on } \Omega^{2h}. \\ \text{Correct fine grid approximation : } v^h \leftarrow v^h + I_{2h}^h e^{2h}. \\ \text{Relax } \nu_2 \text{ times on } A^h u^h = f^h \text{ on } \Omega^h \text{ with initial guess } v^h. \end{array}$ 

Here, in practice, the number of relaxation times  $\nu_1$  and  $\nu_2$  are often 1,2 or 3.

Moreover, the correction scheme is not just two level, it can be more level, by recursive within itself. The algorithm is called the V - cycle. The definition which is given

```
V-cycle Scheme

v^h \leftarrow MV^h(v^h, f^h)

1.Relax \nu_1 times on A^h u^h = f^h with a given initial guess v^h.

2.If \Omega^h = coarsest grid, then go to 4.

Else f^{2h} \leftarrow I_h^{2h}(f^h - A^h v^h).

v^{2h} = 0

v^{2h} \leftarrow MV^{2h}(v^{2h}, f^{2h}).

3.Correct v^h \leftarrow v^h + I_{2h}^h v^{2h}.

4.Relax \nu_2 times on A^h u^h = f^h with initial guess v^h.
```

The V - cycle is just one of a family of multigrid cycling schemes. The entire family

is called the  $\mu - cycle$  method and is defined by

 $\begin{array}{l} \mu\text{-cycle Scheme} \\ \nu^h \leftarrow M\mu^h(v^h, f^h) \\ 1.\text{Relax } \nu_1 \text{ times on } A^h u^h = f^h \text{ with a given initial guess } v^h. \\ 2.\text{If } \Omega^h = \text{coarsest grid, then go to } 4. \\ \text{Else } f^{2h} \leftarrow I_h^{2h}(f^h - A^h v^h). \\ v^{2h} = 0 \\ v^{2h} \leftarrow M\mu^{2h}(v^{2h}, f^{2h}) \ \mu \text{ times.} \\ 3.\text{Correct } v^h \leftarrow v^h + I_{2h}^h v^{2h}. \\ 4.\text{Relax } \nu_2 \text{ times on } A^h u^h = f^h \text{ with initial guess } v^h. \end{array}$ 

If  $\mu = 1$ , which gives the V - cycle and  $\mu = 2$  is called the W - cycle.

Here are other ideas, if it starts on the coarsest discretization with an exact solver, the results are interpolated to the next finer grid with a few cycles (V or W) of the multigrid method are applied .The result is again interpolated to the next finer grid, where again a few cycles of multigrid method are applied.If this is used recursively, the so-called full - multigrid method.The algorithm with V - cycle is given by

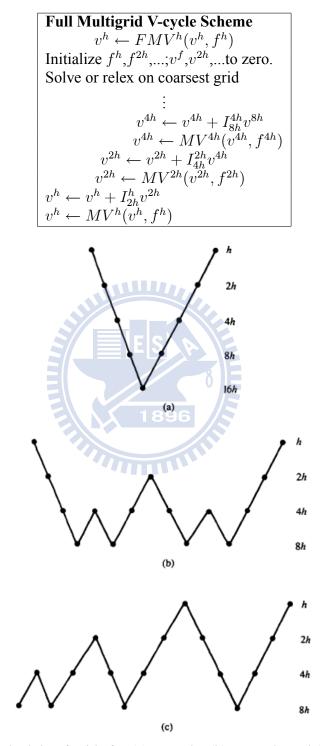


Figure 2.4.1: Schedule of grids for (a) V-cycle, (b) W-cycle and (c) FMV scheme

## 2.5 Complexity

There are various method to solve the linear systems, including stationary iterative methods or nonstationary iterative methods , the following table gives us the complexity of various method :

Table 3.3						
Comple	Complexity of Various Methods					
Method	2D	3D				
JM	$O(N^2)$	$O(N^{\frac{5}{3}})$				
GS	$O(N^2)$	$O(N^{\frac{5}{3}})$				
SOR	$O(N^{\frac{3}{2}})$	$O(N^{\frac{4}{3}})$				
CG	$O(N^{\frac{3}{2}})$	$O(N^{\frac{4}{3}})$				
Multigrid	$O(N \log N)$	$O(N \log N)$				

### 2.6 Numerical Experiments

In this experiments we consider the weighted Jacobi method with  $\omega = 2/3$  applied to the one-dimensional problem Au = 0 on a grid N=48. We use an initial guess,

Experiment 1

$$v_j^h = \frac{1}{2} \left[ \sin(\frac{12j\pi}{N}) + \sin(\frac{30j\pi}{N}) \right]$$

the results of this calculation are given in Figs 2.6.1(a)-(e). The initial guess is shown in Fig 2.6.1(a) . In Fig 2.6.1(b), the approximation  $v^h$  after one relaxation sweep is superimposed on the initial guess. Much of the oscillatory component of the initial guess has already been removed. The maximum norm of the error has decreased signification. Fig. 2.6.1(c) shows the approximation after three relaxation sweeps, superimposed on the previous approximations. Further relaxations on the fine grid would provide only a slow improvement at this

point. This signals that it is time to move to the coarse grid.Fig. 2.6.1(d) shows the fine grid error after one relaxation sweep on the coase grid and the error after three coarse grid relaxation sweeps is shown in Fig. 2.6.1(e)

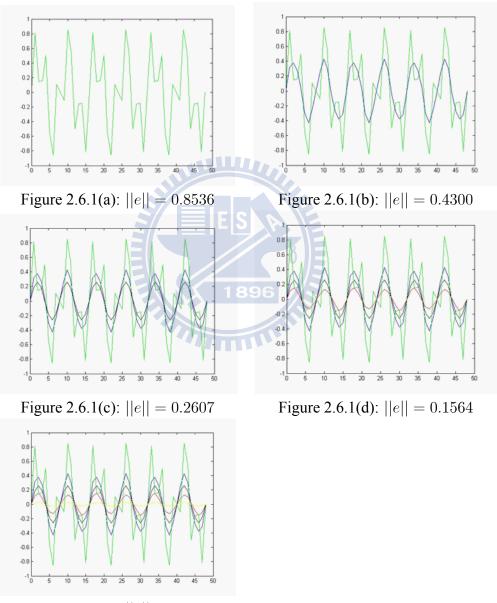


Figure 2.6.1(e): ||e|| = 0.0243

#### **Experiments 2**

Consider the following two-point boundary value problem. It is given by the secondorder ordinary differential equation

$$\begin{cases} (\alpha(2+\cos(\beta x))u'(x))' = f(x), & 0 < x < \pi, \ \alpha \text{ and } \beta \text{ are constant.} \\ u(0) = 0, \ u(\pi) = \pi^2 \end{cases}$$
(2.83)

we use finite difference method to discrete the domain. The domain of the problem  $\{x: 0 \le x \le \pi\}$  is partitioned into N subintervals by introducing the grid points  $x_j = jh$ , where  $h = \pi/N$  is the constant width of the subintervals. At each of the N - 1 interior grid points, the differential equation is replaced by a second-order finite difference approximation. We also introduce  $v_j$  as an approximation to the exact solution  $u(x_j)$ .

$$\frac{\left[2\alpha + \alpha \cos(\beta x_{j-\frac{1}{2}})]v_{j-1} - \left[4\alpha + \alpha \cos(\beta x_{j+\frac{1}{2}}) + \alpha \cos(\beta x_{j-\frac{1}{2}})\right]v_j}{h^2} + \frac{h^2}{1 + \alpha \cos(\beta x_{j+\frac{1}{2}})]v_{j+1}} = f(x_j), \quad 1 \le j \le N-1.$$

$$v_0 = 0, \ v_N = \pi^2$$
(2.84)

Let a represent the entry  $4\alpha + \alpha \cos(\beta x_{j+\frac{1}{2}}) + \alpha \cos(\beta x_{j-\frac{1}{2}})$ , b represent the entry  $2\alpha + \alpha \cos(\beta x_{j+\frac{1}{2}})$ , c represent the entry  $2\alpha + \alpha \cos(\beta x_{j-\frac{1}{2}})$ , the system can be represent in matrix form as Av = f.

$\frac{1}{h^2}$	$\begin{bmatrix} a \\ c \end{bmatrix}$	$b \\ a \\ \cdot .$	b •	•.		$ \begin{bmatrix} v_1 \\ \vdots \\ \vdots \end{bmatrix} $		$\begin{array}{c} f_1 \\ \vdots \\ \vdots \end{array}$	
h²			С	$a \\ c$	$\begin{bmatrix} b\\ a \end{bmatrix}$	$v_{N-1}$		$\vdots$ $f_{N-1}$	

The matrix A is tridagonal, symmetric positive definite and has dimension  $(N-1) \times (N-1)$ .

We use multigrid method to solve the linear system. There are several comments in order. First, we choose  $\alpha$  and  $\beta$  the constant one. Second, we use three level v - cycle and

the number of relaxation times  $\nu_1$  and  $\nu_2$  is the constant three. Third, we use  $v = \pi x$  as an initial guess.

The solution of this boundary value problem is  $u(x) = x^2$ . The error  $e^h$  is defined by  $e^h = ||u^h - v^h||_{L^2}$ , where  $u^h$  is the exact solution and  $v^h$  is the approximation on  $\Omega^h$ . The rate of convergence  $\sigma$  is computed by

$$\sigma = \frac{e^{2h}}{e^h} \tag{2.85}$$

Table2.6.1					
Ν	e	σ			
128	2.7315e-004				
256	6.8301e-005	3.9992			
512	1.7076e-005	3.9998			

The results are tabulated in above .The theoretical convergence rate is 4 for  $\sigma$ .

#### 1896 Experiments 3

by

In experiment two, we consider the boundary value problem in two-dimension given

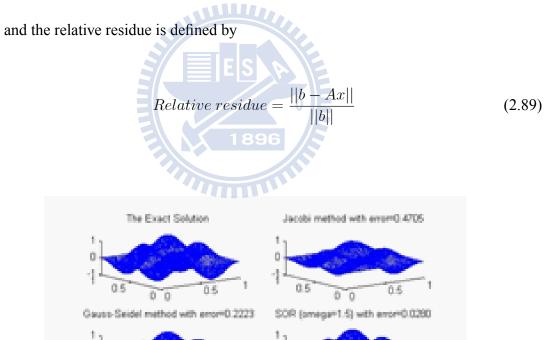
$$-\bigtriangleup u = f \quad \text{in} \quad \Omega : [0,1] \times [0,1]$$
$$u = 0 \quad \text{on boundary}$$
(2.86)

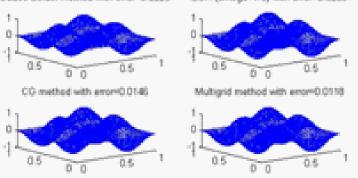
and finite element method is used to discretize the domain. To solve the linear systems Ax = b, we applied the Jacobi Method, Gauss-Seidel Method, Successive Overrelaxation Method, Conjugate Gradient Method and Multigrid Method. Compare the iteration numbers and the relative residual error between the multigrid method and other methods , we can find the advantages of using multigrid method as the linear systems solver. The exact solution is

$$u(x,y) = \sin(2\pi x)\sin(4\pi y)$$
 (2.87)

The tolerance is 10e-16. The error is defined by

$$error = ||u_{ex} - \hat{u}||_{\infty} \tag{2.88}$$





Figrue2.6.2: Approximate solutions

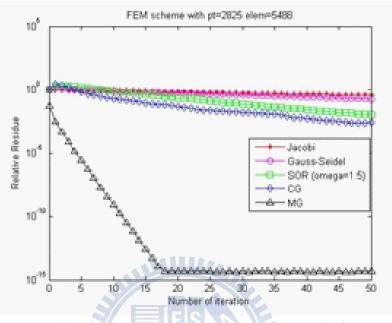


Figure 3.6.1 show the exact solution and five kinds of approximation.

Figure 2.6.3: Iteration numbers of five methods

Figure 3.6.2 show us the relation between the number of iteration and the log relative

residue. With five kinds of iteration method, Multigrid Method gets the best performance.

# Chapter 3 Research Method

Let  $\Omega$  be a bounded polygonal domain in  $\mathbb{R}^2$  with re-entrant angle. Consider the Poisson equation with homogeneous boundary condition

$$\begin{cases} -\Delta u = f & \text{in } \Omega\\ u = 0 & \text{on } \partial \Omega \end{cases}$$
(3.90)

where  $f \in L^2(\Omega)$ .

When (3.90) is solved by the P<sub>1</sub> finite element method on a quasi-uniform grid ,the convergence rate in the energy norm is therefore of order  $O(h^{(\pi/\omega)-\epsilon})$ , where h is the mesh size of the triangulation and  $\omega$  is the re-entrant angle.

In 1996, S.C.Brenner improves the convergence rate developed in [3]. It is based on the full multigrid iteration technique and the following singular function representation of u [4, 5, 6, 7]

$$u = \kappa s + w \tag{3.91}$$

where  $w \in H^2(\Omega)$  and the *s* are the singular function associated with the re-entrant angle. Note that the coefficient  $\kappa$  is known as stress intensity factors in elasticity problems.

The multigrid method in [3] computes a solution of (3.91) in the form of

$$u_h = \kappa_h s + w_h$$

where  $w_h$  is a piecewise linear function. It is shown in [3] that

$$|u - u_h|_{H^1(\Omega)} \le Ch||f||_{L^2(\Omega)},\tag{3.92}$$

$$|\kappa - \kappa_h| \le Ch^{1 + \pi/\omega - \epsilon} ||f||_{L^2(\Omega)}.$$
(3.93)

In 1997, S.C.Brenner and L.-Y.Sung [9] extend the results in [3] and [10] to the case of a polygonal domain with crack in  $\mathbb{R}^2$  (i.e. the re-entrant angle  $\omega = 2\pi$ .)

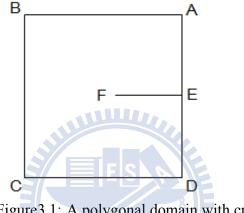


Figure 3.1: A polygonal domain with cracks.

Let p be the vertice of  $\Omega$  such that the angle  $\omega$  associated with p satisfies  $\omega > \pi$  (i.e. the vertice F in the figure 1.1). Let polar coordinates  $(r, \theta)$  be chosen at the vertex p so that the angle  $\omega$  is spanned by the two rays  $\theta = 0$  and  $\theta = \omega$ .

The singular function s is defined by

$$s(r,\theta) = \phi(r)r^{\pi/\omega}\sin(\frac{\pi}{\omega}\theta)$$
(3.94)

where  $\phi(r)$  is smooth cut-off function which equal 1 identically in a neighborhood of 0, and the supports of the  $\phi$  is small enough so that the singular function s vanish identically on  $\partial \Omega$ .

Then the solution u has the representation

$$u = \kappa s + w \tag{3.95}$$

The stress intensity fractors  $\kappa$  can be expressed in terms of u by the following extraction formula

$$\kappa = \frac{1}{\pi} \left\{ \int_{\Omega} f s_{-1} dx + \int_{\Omega} u \bigtriangleup s_{-1} dx \right\},\tag{3.96}$$

where the dual singular function  $s_{-1}$  is defined in the polar coordinate system  $(r, \theta)$  as

$$s_{-1}(r,\theta) = \phi(r)r^{-\pi/\omega}\sin(\frac{\pi}{\omega}\theta)$$

([5, 7, 11, 12, 13, 14, 15, 16]).

There is a idea that we will take advantage of the singular function representation. We substitute the representation (3.95) into (3.90) to obtain the following boundary-value problem for w:

$$\begin{cases} -\Delta w = f + \kappa \Delta s & \text{in } \Omega \\ w = 0 & \text{on } \partial \Omega \end{cases}$$
(3.97)

If the  $\kappa$  were known, we could solve (3.97) using piecewise linear finite element method. Unfortunately the  $\kappa$  is unknown, so we apply the finite element method on the kth level to the following varational problem:

Find  $w_k \in H_0^1(\Omega)$  such that

$$\int_{\Omega} \nabla \widehat{w}_k \nabla v dx = \int_{\Omega} \left( f + \kappa_k \bigtriangleup s \right) v dx \ \forall v \in H^1_0(\Omega)$$
(3.98)

where the approximate stress intensity factors  $\kappa_k$  is computed by the extraction formula (3.96) using the approximate solution  $u_{k-1}$  obtained in the (k-1)st level, i.e.,

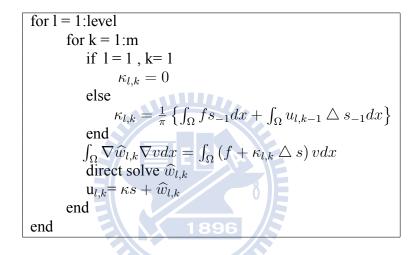
$$\kappa_k = \frac{1}{\pi} \left\{ \int_{\Omega} fs_{-1} dx + \int_{\Omega} u_{k-1} \bigtriangleup s_{-1} dx \right\}.$$
(3.99)

We obtain, on the kth level, a piecewise linear approximate solution  $w_k$  to  $\hat{w}_k$  by applying the kth level iteration n times using  $w_{k-1}$  as the initial guess. The approximate solution  $u_k$  to u is defined to be

$$u_k = \kappa_k s + w_k. \tag{3.100}$$

In other words we are really computing the regular part w of the solution. The improvement in the convergence rate is possible because w has better regularity than u.

The algorithm of S.C.Brenner's method is given by the following



Although we known that the singular function can replace the solution at the sigular point well, but there still have some problem we can observe. After leaving the singular point, the singular function is now not a correct solution ,so the error occured on this cutoff region. To improve this problem, we introduce two strategys, adaptive mesh-refinement techniques and adaptive cut-off function.

For the first strategy, adaptive mesh-refinement techniques, there is a relationship between the converge rate and the error estimator, we should find suitable order of h for the error estimator. The converge rate in the ragular domain is

$$||u - u_h||_{L^2(\Omega)} \leq Ch^2 |u|_{H^2(\Omega)}$$
  
$$||u - u_h||_{H^1(\Omega)} \leq Ch |u|_{H^2(\Omega)}, \qquad (3.101)$$

the order of error estimator is choose by

$$E(K) = \alpha ||h(f - au)||_{K} + \beta (\frac{1}{2} \sum_{\tau \in \partial K} h_{\tau} [n_{\tau} \cdot (c\nabla u_{h})]^{2})^{1/2}$$
(3.102)

Now, in the sigular domain , we also want to find the same relationship between the converge rate and the error estimator. Since the converge rate in singular domain is

$$||u - u_h||_{L^2(\Omega)} \leq Ch^{(3/2) - \epsilon} |u|_{H^2(\Omega)}$$
  
$$||u - u_h||_{H^1(\Omega)} \leq Ch^{(1/2) - \epsilon} |u|_{H^2(\Omega)}, \qquad (3.103)$$

the order of error estmator is choose by

$$E(K) = \alpha ||h^{(3/4)}(f - au)||_{K} + \beta (\frac{1}{2} \sum_{\tau \in \partial K} h_{\tau}^{(1/4)} [n_{\tau} \cdot (c\nabla u_{h})]^{2})^{1/2}$$
(3.104)

we use this order as the rules of adaptive mesh-refinement techniques.

For the second strategy, adaptive cut-off function .Since we realize that the singular function replaced the solution successfully only near the singular point, so the range of cut-off function may affect the error. If we choose the range of cut-off function is too wide, most solution replace by the singular function , that is not a correct .Otherwise, if we choose the range of cut-off functions is too small, the error still big near the singular point.So, this adaptive cut-off function strategy is find the suitable cut range and move forward slowly .On the one hand, the error will not so big . On the other hand, the solution replace by the singular function with respect to the mesh-refinement steps.

The adaptive cut-off funciton is defined by

$$\phi(r) = a_0 + a_1 r + a_2 r^2 + a_3 r^3 + a_4 r^4 + a_5 r^5 \tag{3.105}$$

find the coefficient  $a_0, a_1, a_2, a_3, a_4, a_5$  which is satisfy

4.

$$\begin{cases} \phi(\frac{1}{3+i}) = 1 , & \phi(\frac{2}{3+i}) = 0 \\ \phi'(\frac{1}{3+i}) = 0 , & \phi'(\frac{2}{3+i}) = 0 \\ \phi''(\frac{1}{3+i}) = 0 , & \phi''(\frac{2}{3+i}) = 0 \end{cases}$$
(3.106)

where  $\frac{1}{3+i} \le r \le \frac{2}{3+i}, i = 0, 1, 2$  is the mesh-refinement steps.

The purpose of this range we choose is that we want the interval of cut range is reduce and close to singular point, also the cut range is move forward slowly.

From the two strategy above, we can hold the stress intensity factorys and also improve the accuracy of the global error. The results shown in the experiment 4 in chapter



# **Chapter 4 Numerical Results**

In this chapter we report the result of some numerical experiments. The first experiment we show the error arise from the singular point associated with the re-entrant angle.We consider the following boundary value problem:

$$\begin{cases} -\Delta u = f & \text{in } \Omega\\ u = 0 & \text{on } \partial \Omega \end{cases}$$
(4.107)

where  $\Omega$  is the circle with four angle  $\omega = \pi/2, 3\pi/2, 7\pi/4$  and  $\pi/0.51, \omega$  is the maximum re-entrant angle.

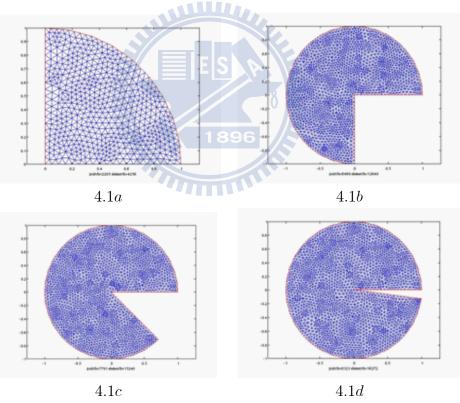


Figure 4.1: Circle with four kinds angle.

Let the exact solution u be

$$u = (1 - r^2)r^\beta \sin(\beta\theta) \tag{4.108}$$

where  $0 < r < 1, 0 < \theta < \omega, \beta = \frac{\pi}{\omega}, \omega$  is the maximum re-entrant angle. When is solved by piecewise linear finite element method on a quasi-uniform grid, the error is shown in the following:

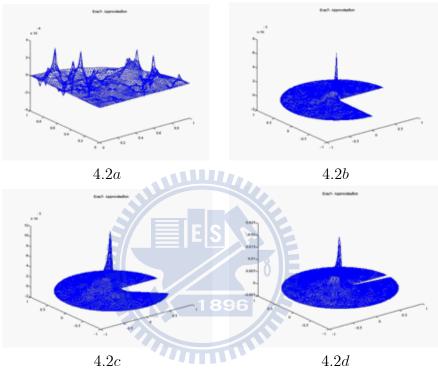


Figure 4.2: Error with four kinds of angle

Table 4.1 Results for four kinds of angle						
$\pi$ $3\pi$ $7\pi$ $\pi$						
	$\omega = \frac{1}{2}$	$\omega = -\frac{1}{2}$	$\omega = -4$	$\omega = \frac{1}{0.51}$		
$  u-u_h  _{H^1(\Omega)}$	2.6377E-02	4.1243E-02	7.0759E-02	9.6003E-02		
# of points	2205	6469	7791	8323		

we can see the Fig4.2a is not a re-entrant angle, so the error of Fig4.2a didn't have peak at the origin, but the others have. We also can find when the maximum re-entrant angle  $\omega$  arise, the error also arise with the  $\omega$ .

The second experiment we will concern about the mesh-refinement strategy. There are three different mesh-refinement strategy we used. The first strategy is uniform mesh

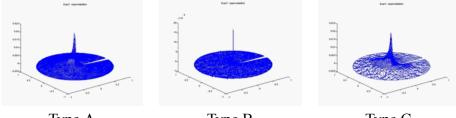
with uniformly refinement, the second is singular element mesh with uniformly refinement and the last is uniform mesh with adaptive short cut-region refinement.

The first mesh-refinement strategy is like as we refine mesh as usual. The second strategy, we changed the uniform mesh to the singular mesh. According to our mentioned in the singular element, here the approach is like singular element, we put the exponential grid points (x, y) in the way

$$\begin{cases} (x,y) = (2^{-i},0) \\ (x,y) = (2^{-i}\cos\omega, 2^{-i}\sin\omega) \end{cases}, \ i = 1,...,10$$
(4.109)

on the two rays. This may couse many grid points located near the singular point. The third strategy is using uniform mesh, but not refine the whole domain, only refine the region with radius  $r = 0.5^{j}$  form the center points (0,0), where j represent the times of refinement.

We consider the maximum re-entrant angle  $\omega = \frac{\pi}{0.51}$  case with three different meshrefinement strategys .The times of mesh-refinement steps is 2, means that we have the original corasest domain  $\Omega^{4h}$ , find domain  $\Omega^{2h}$  and  $\Omega^{h}$ .



Type AType BType CFigure 4.3: Error of three types of mesh-refinement

we caculate the H1-norm error and L2-norm in the following table

Table4.1					
Type A Type B Type C					
$  u - u_h  _{H^1(\Omega)}$	9.6003E-02	2.9259E-02	1.3185E-01		
$  u-u_h  _{L^2(\Omega)}$	3.1643E-03	2.3718E-04	4.9345E-03		
# of points	8323	15085	1373		

From the Table4.1, we discuss some phenomena. In the second strategy, there are many points around of center. When we refine the mesh, the points around the center of all will be refined. Although we reduced the error, but we also pay a high price since too many points cause the large matrix systems. In the third strategy, we only refine the points inside the cut-region, the error outside the cut-region still not be reduced, therefore, the accuracy performance is not good. The convergence rate for type A and type B in the  $H^1 - norm$  is therefore of order  $O(h^{(\pi/\omega)-\epsilon}).(i.e.$  the theoratical convergence rate is 1.42)

	Table4.2						
Ω	Type A	ratio A	Type B	ratio B	Type C	ratio C	
$\Omega^{4h}$	2.3184E-01		1.1065E-01		2.1384E-01		
$\Omega^{2h}$	1.4147E-01	1.5116	5.6813E-02	1.9476	1.6135E-01	1.3253	
$\Omega^h$	9.6003E-02	1.4737	2.9259E-02	1.9417	1.3185E-01	1.2237	

Simultaneously, we apply the multigrid method for solving the linear systems and observed the benefits for three different mesh-refinement. For the multigrid parameter, we apply V-cycle 2th level iteration and we choose the weight-Jacobi method to relaxation, the maximum relaxation number is 2.For the restriction operator ,we use full weighting as a restriction operator, and for the interpolation operator, we consider linear interpolation.

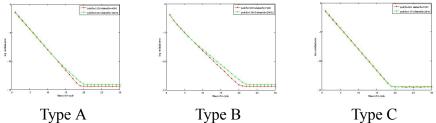


Figure 4.4: Multigrid benefits for three different mesh-refinement

From the Figure 4.4 shown in above, we can observe that although the mesh size is different, the slope almost not changed, meens that the converge rate of multigrid method is indenpentent of mesh size. The third experiment we use the method with S.C.Brenner and L.-Y.Sung.The singular function for  $\Omega$  is

$$s(r,\theta) = \phi(r)r^{\pi/0.51}\sin(\frac{\pi}{0.51}\theta)$$
 (4.110)

where the cut-off function  $\phi(r)$  is defined to be

$$\begin{cases} 1, 0 \le r \le \frac{1}{3} \\ -1458r^5 + 3645r^4 - 3510r^3 + 1620r^2 - 360r + 32, \frac{1}{3} \le r \le \frac{2}{3} \\ 0, \frac{2}{3} \le r \end{cases}$$
(4.111)

The  $\kappa_k$  are obtained by the extraction formula

$$\kappa_k = \frac{1}{\pi} \left\{ \int_{\Omega} fs_{-1} dx + \int_{\Omega} u_{k-1} \bigtriangleup s_{-1} dx \right\}.$$
(4.112)

where

$$s_{-1}(r,\theta) = \phi(r)r^{-\pi/0.51}\sin(\frac{\pi}{0.51}\theta).$$
(4.113)

Here, for the approximate stress intensity factors  $\kappa_k$ , we choose 3-th level (l = 3) and itera-

tion 5 times (m = 5)

The error is shown in the following

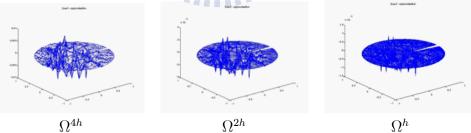


Figure 4.5: Error of three level mesh

The theoretical number for the stress intensity factorys  $\kappa$  is 1.

	Table4.3					
Ω	$\kappa$	$   u-u_h  _{H^1(\Omega)}$	# of points			
$\Omega^{4h}$	9.9399E-01	2.3664E-01	556			
$\Omega^{2h}$	9.9815E-01	1.5423E-01	2128			
$\Omega^h$	9.9944E-01	1.0451E-01	8323			

The fourth experiment we improve the method by including adaptive mesh-refinement

techniques and adaptive cut-off function. The following table shown the stress intensity fac-

	Table4.4				
Ω	$\kappa$	$  u-u_h  _{H^1(\Omega)}$	# of points		
1	9.9399E-01	2.3664E-01	556		
2	1.0017E+01	2.0107E-01	913		
3	9.9098E-01	1.7432E-01	1359		
4	9.8648E-01	1.4946E-01	1842		
5	9.8810E-01	1.3034E-01	2196		
6	9.8722E-01	1.2721E-01	2747		
7	9.8841E-01	1.1381E-01	3207		
8	9.9204E-01	1.0953E-01	3533		
9	9.9492E-01	1.0199E-01	4522		
10	9.9341E-01	9.8286E-02	5494		
11	9.9306E-01	9.7071E-02	6178		
12	9.9560E-01	9.5075E-02	7217		
13	9.9534E-01	9.5117E-02	7708		
14	9.9596E-01	9.3322E-02	9533		
15	9.9626E-01	9.3082E-02	10568		

torys  $\kappa$  and the error in  $H^1 - norm$ .

we can clearly compare the  $||u - u_h||_{H^1(\Omega)}$  with these two results above. Fixed points on

both sides of almost equal, we can clearly see the improvement in error in  $H^1 - norm$ .

	Table4.5						
	S.C.Brenn	Experimer	nt				
Ω	$\Omega \mid   u - u_h  _{H^1(\Omega)} \mid \# \text{ of points}$			$  u - u_h  _{H^1(\Omega)}$	# of points		
1	2.3664E-01	556	1	2.3664E-01	556		
2	1.5423E-01	2128	4	1.4946E-01	1842		
3	1.0451E-01	8323	13	9.5117E-02	7708		

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