

## Part 2: Multigrid Methods for the computation of singular solutions and stress intensity factors

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## Used literature



S. C. BRENNER, *Multigrid Methods for the computation of singular solutions and stress intensity factors : Corner singularities I.*, Department of Mathematics and Center for Computation and Technology, Mathematics of computation, April 1999, Volume 86, Number 226, Pages 559-583.

# Introduction

Let be:

$\{\mathcal{T}_k\}$ ,  $k \geq 1$ , a family of triangulations of  $\Omega$ , where a regular subdivision  $\mathcal{T}_{k+1}$  of is obtained from  $\mathcal{T}_k$  by connecting the edges of the triangles in  $\mathcal{T}_k$ .

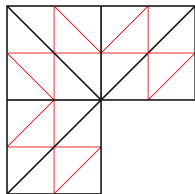


Figure: Triangulation

$V_k = \{v \in H_0^1(\Omega) : v|_T \in \mathcal{P}_1 \ \forall T \in \mathcal{T}_k\}$  ... piecewise linear finite elements associated with  $\mathcal{T}_k$ .

# Introduction

The **discrete inner product**  $(\cdot, \cdot)_k$  defined by

$$(v_1, v_2)_k = h_k^2 \sum_{\text{vertices } p \text{ of } \mathcal{T}_k} v_1(p)v_2(p) \quad \forall v_1, v_2 \in V_k.$$

$\Rightarrow (v, v)_k$  is spectral equivalent to  $\|v\|_{L^2(\Omega)}^2 \quad \forall v \in V_k.$

The **operators**  $A_k : V_k \rightarrow V_k$  and  $I_k^{k-1} : V_k \rightarrow V_{k-1}$  (Restriction Operator), defined by:

$$(A_k v_1, v_2)_k = \int_{\Omega} \nabla v_1 \cdot \nabla v_2 \, dx \quad \forall v_1, v_2 \in V_k \subset V_{k-1},$$

$$(I_k^{k-1} v, w)_{k-1} = (v, w)_k \quad \forall v \in V_k, w \in V_{k-1}.$$

$\Rightarrow A_k$  symmetric, positive definite and the spectralradius  $\rho(A_k) \lesssim h_k^{-2}.$

# Standard $k$ -th level multigrid iteration

The  $k$ -th level multigrid iteration with initial guess  $z_0$  yields  $MG(k, z_0, g)$  as an approximate solution to the equation

$$A_k z = g.$$

For  $k = 1$ ,  $MG(1, z_0, g)$  is the solution obtained from an exact solver, i.e.  $MG(1, z_0, g) = A_1^{-1}g$ .

For  $k > 1$ , there are two steps.

**Smoothing Step:** Let  $z_l \in V_k$  ( $1 \leq l \leq m$ ) be defined recursively by the equations

$$z_l = z_{l-1} + \frac{1}{\gamma_k} (g - A_k z_{l-1}), \quad 1 \leq l \leq m, \quad (\text{Richardson Relaxation})$$

where  $m \in \mathbb{N}_0$  independent of  $k$ , and  $\gamma_k = Ch_k^{-2}$  dominates  $\rho(A_k)$ .

**Correction Step:** Let  $\bar{g} = I_k^{k-1}(g - A_k z_m) \in V_{k-1}$  and  $q_i \in V_{k-1}$  ( $0 \leq i \leq p$ ,  $p = 1$  (V-cycle) or  $p = 2$  (W-cycle)) be defined recursively by

$$q_0 = 0 \quad \text{and} \quad q_i = \underbrace{MG(k-1, q_{i-1}, \bar{g})}_{\text{approx. of } A_{k-1}^{-1} \bar{g}}, \quad 1 \leq i \leq p.$$

The output is obtained by combining the two steps:

$$MG(k, z_0, g) = z_m + q_p.$$

# Full multigrid algorithm 1

If  $f \in L^2(\Omega)$ , we use the nested iteration to compute  $\kappa_k$  and  $w_k$ .

The nested iteration:

For  $k = 1$ ,

$$w_1 = A_1^{-1} g_1, \quad \text{where } (g_1, v)_1 = \int_{\Omega} f v \, dx \quad \forall v \in V_1.$$

We set

$$\kappa_1 = 0 \quad \text{and} \quad u_1 = w_1.$$

For  $k \geq 2$ ,  $\kappa_k \in \mathbb{R}$  are computed by

$$\kappa_k = \frac{1}{\pi} \left( \int_{\Omega} f s_- \, dx + \int_{\Omega} u_{k-1} \Delta s_- \, dx \right),$$

and  $w_k \in V_k$  is obtained recursively by

$$w_{k,0} = w_{k-1}, \quad w_{k,l} = MG(k, w_{k,l-1}, g_k), \quad w_k = w_{k,n} \quad \text{for } 1 \leq l \leq n,$$

where  $n$  is a positive integer independent of  $k$ , and  $g_k \in V_k$  is defined by

$$(g_k, v)_k = \int_{\Omega} (f v \, dx + \kappa_k \Delta s_+) \, dx \quad \forall v \in V_k.$$

We define then  $u_k$  by

$$u_k = \kappa_k s_+ + w_k.$$

# Full multigrid algorithm 2

If  $f \in H^1(\Omega)$ , we use the nested iteration to compute  $\kappa_{\ell,k}$ ,  $\ell \in \mathcal{L}$  and  $w_k$ .

The **nested iteration**:

For  $k = 1$ ,

$$w_1 = A_1^{-1} g_1, \quad \text{where} \quad (g_1, v)_1 = \int_{\Omega} f v \, dx \quad \forall v \in V_1.$$

We set

$$\kappa_{\ell,1} = 0 \quad \text{for } \ell \in \mathcal{L}, \quad \text{and} \quad u_1 = w_1.$$

For  $k \geq 2$ ,  $\kappa_{\ell,k} \in \mathbb{R}$  are computed by

$$\kappa_k = \frac{1}{\ell\pi} \left( \int_{\Omega} f s_{+,-\ell} \, dx + \int_{\Omega} u_{k-1} \Delta s_{+,-\ell} \, dx \right) \quad \text{for } \ell \in \mathcal{L},$$

and  $w_k \in V_k$  is obtained recursively by

$$w_{k,0} = \mathcal{J}_{k-1}^k w_{k-1}, \quad w_{k,l} = MG(k, w_{k,l-1}, g_k), \quad w_k = w_{k,n} \quad \text{for } 1 \leq l \leq n,$$

where  $n$  is a positive integer independent of  $k$ , and  $g_k \in V_k$  is defined by

$$(g_k, v)_k = \int_{\Omega} \left( f v \, dx + \sum_{\ell \in \mathcal{L}} \kappa_{\ell,k} \Delta s_{+,\ell} \right) \, dx \quad \forall v \in V_k.$$

We define then  $u_k$  by

$$u_k = \sum_{\ell \in \mathcal{L}} \kappa_{\ell,k} s_{+,\ell} + w_k.$$

# The intergrid transfer operator $\mathcal{J}_{k-1}^k$

For

$$Q_k \subset H_0^1(\Omega) \quad (k = 0, 1, 2, \dots) \text{ quadratic Lagrange finite element space associated with } \mathcal{T}_k. \quad (1)$$

we define the interpolation operators

$$\begin{aligned} \mathcal{I}_{k-2}^{k-1} : Q_{k-2} &\rightarrow V_{k-1} \\ w &\rightarrow v, \text{ such that } v(p) = w(p) \quad \forall \text{ vertices } p \text{ of } \mathcal{T}_{k-1} \end{aligned}$$

which is an isomorphism and

$$\begin{aligned} \mathcal{I}_{k-2}^k : Q_{k-2} &\rightarrow V_k \\ w &\rightarrow v, \text{ such that } v(p) = w(p) \quad \forall \text{ vertices } p \text{ of } \mathcal{T}_k, \end{aligned}$$

and further the **intergrid transfer operator**:

$$\mathcal{J}_{k-1}^k = \mathcal{I}_{k-2}^k \circ (\mathcal{I}_{k-2}^{k-1})^{-1} : V_{k-1} \rightarrow V_k \quad \text{for } k = 2, 3, \dots$$



# Contraction properties for the $k$ -th level iteration

Convergence result for the  $k$ -th level iteration in the energy norm:

## Lemma

Let  $p = 1$  (V-cycle) or  $p = 2$  (W-cycle) and  $m \geq 1$  in the  $k$ -th level iteration. Then there exists a  $\delta < 1$ , independent of  $k$ , such that

$$|z - MG(k, z_0, g)|_{H^1(\Omega)} \leq \delta |z - z_0|_{H^1(\Omega)}. \quad (2)$$

Convergence result for the  $k$ -th level iteration in the  $\|\cdot\|_{H^{1-(\pi/\omega)+\epsilon}(\Omega)}$  norm:

## Theorem

Let  $p = 2$  (W-cycle),  $0 < \delta < 1$ ,  $0 < \epsilon < \pi/\omega$  and  $\alpha_\epsilon = 1 - \pi/\omega \neq 1/2$ . If the number of smoothing steps  $m$  in the  $k$ -th level iteration is sufficiently large, then we have

$$\|z - MG(k, z_0, g)\|_{H^{\alpha_\epsilon}(\Omega)} \leq \delta \|z - z_0\|_{H^{\alpha_\epsilon}(\Omega)}. \quad (3)$$

# Convergence Analysis for the full multigrid algorithm 1

## Theorem

Let  $p = 2$  (W-cycle),  $0 < \epsilon < \pi/\omega$ ,  $\alpha_\epsilon = 1 - \pi/\omega \neq 1/2$  and the number of smoothing steps  $m$  in the  $k$ -th level iteration be sufficiently large, that (2) and (3), hold for  $0 < \delta < 1$ . If the number of nested iterations  $n$  is sufficiently large, then we have

$$|w - w_k|_{H^1(\Omega)} \lesssim h_k \|f\|_{L^2(\Omega)}, \quad (4)$$

$$|\kappa - \kappa_k| \lesssim_\epsilon h_k^{1+\pi/\omega-\epsilon} \|f\|_{L^2(\Omega)}, \quad (5)$$

$$\|w - w_k\|_{H^{\alpha_\epsilon}(\Omega)} \lesssim_\epsilon h_k^{1+\pi/\omega-\epsilon} \|f\|_{L^2(\Omega)}. \quad (6)$$

where  $w_k$  and  $\kappa_k$  are computed by **Full multigrid algorithm 1**.

## Corollary

Under the assumption of theorem 3, we have

$$|u - u_k|_{H^1(\Omega)} \lesssim h_k \|f\|_{L^2(\Omega)}. \quad (7)$$

# Convergence Analysis for the full multigrid algorithm 2

Properties of the intergrid transfer operator  $\mathcal{J}_{k-1}^k : V_{k-1} \rightarrow V_k$

## Lemma

We have the following estimates for  $\mathcal{J}_{k-1}^k$ :

$$|\mathcal{J}_{k-1}^k v|_{H^1(\Omega)} \lesssim h_k |v|_{H^1(\Omega)} \quad \forall v \in V_{k-1}, \quad (8)$$

$$|\Pi_k \eta - \mathcal{J}_{k-1}^k \Pi_{k-1} \eta|_{H^1(\Omega)} \lesssim h_k^{1+t} \|\eta\|_{H^{2+t}(\Omega)} \quad \forall \eta \in V_{k-1}, \quad (9)$$

where

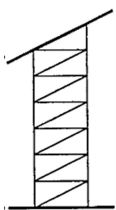
$$\begin{aligned} \Pi_k : H^1(\Omega) &\rightarrow V_k \\ w &\rightarrow v, \text{ such that } v(p) = w(p) \quad \forall \text{ vertices } p \text{ of } \mathcal{T}_k, \end{aligned}$$

is the nodal interpolation operator associated with  $V_k$  and  $0 \leq t \leq 1$ .

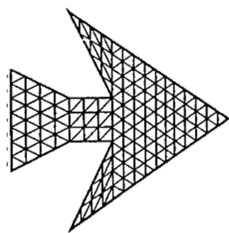
# Uniform band condition (UBC)

## Definition

A **uniform band** in a triangulation is a collection of triangles between two parallel lines, such that any two triangles sharing a common side form a parallelogram (see figure 2). We say a triangulation satisfies the **uniform band condition (UBC)**, if it can be divided completely into uniform bands (see figure 3).



Figure



Figure

## Remark

One can always find a triangulation satisfying the uniform band condition for any **polygonal domain** whose vertices all have rational coordinates, and the uniform band condition is preserved by **regular subdivision**.

# Super convergence result

We define the **Ritz projection operator**  $P_k : H^2(\Omega) \rightarrow V_k$  by

$$\int_{\Omega} \nabla(\eta - P_k \eta) \cdot \nabla v \, dx = 0 \quad \forall \eta \in H_0^1(\Omega), v \in V_k.$$

## Lemma

*Suppose the triangulations  $\mathcal{T}_k$  satisfy the uniform band condition and  $\eta \in H^3(\Omega) \cap H_0^1(\Omega)$ . Then*

$$|\Pi_k \eta - P_k \eta|_{H^1(\Omega)} \lesssim h_k^2 \|\eta\|_{H^3(\Omega)}.$$

## Corollary

*Suppose the triangulations  $\mathcal{T}_k$  satisfy the uniform band condition and  $\eta \in H^3(\Omega) \cap H_0^1(\Omega)$  for  $0 \leq t \leq 1$ . Then*

$$|\Pi_k \eta - P_k \eta|_{H^1(\Omega)} \lesssim h_k^{1+t} \|\eta\|_{H^{2+t}(\Omega)}.$$

# Convergence Analysis for the full multigrid algorithm 2

Using this previous **superconvergence result** we can show:

## Theorem

Let  $f \in H^1(\Omega)$ . Assume that the triangulations  $\mathcal{T}_k$  satisfy the uniform band condition,  $p = 1$  (V-cycle), or  $p = 2$  (W-cycle), and  $m \geq 1$ . If the number of nested iterations  $n$  is sufficiently large, then we have

$$|\Pi_k w - w_k|_{H^1(\Omega)} \lesssim_\epsilon h_k^{2-\epsilon} \|f\|_{H^1(\Omega)}, \quad (10)$$

$$\sum_{\ell \in \mathcal{L}} |\kappa_\ell - \kappa_{\ell,k}| \lesssim_\epsilon h_k^{2-\epsilon} \|f\|_{H^1(\Omega)}, \quad (11)$$

where  $\mathcal{L} = \{\ell \in \mathbb{N} : \ell\pi/\omega < 2\}$  and  $w_k, \kappa_{\ell,k}$  are computed by **Full multigrid algorithm 2**.

## Corollary

Under the assumption of theorem 7, we have

$$|u - u_k|_{H^1(\Omega)} \lesssim h_k \|f\|_{H^1(\Omega)} \quad \text{and} \quad \max_p |u(p) - u_k(p)|_{H^1(\Omega)} \lesssim_\epsilon h_k^{2-\epsilon} \|f\|_{H^1(\Omega)}, \quad (12)$$

where the maximum is taken over all the vertices  $p \in \mathcal{T}_k$ .

## Remark

If for all internal angles  $\omega$  of  $\Omega$  we have  $\ell\omega \neq \pi/2$  for all  $\ell \in \mathbb{N}$ , then

$$w \in H^3(\Omega),$$

from which follows the  *$\epsilon$ -independent* estimates:

$$|\Pi_k w - w_k|_{H^1(\Omega)} \lesssim h_k^2 \|f\|_{H^1(\Omega)},$$

$$\sum_{\ell \in \mathcal{L}} |\kappa_\ell - \kappa_{\ell,k}| \lesssim h_k^2 \|f\|_{H^1(\Omega)},$$

$$\max_p |u(p) - u_k(p)|_{H^1(\Omega)} \lesssim h_k^2 |\ln h_k|^{1/2} \|f\|_{H^1(\Omega)}.$$

# Model data:

Domain  $\Omega$ :

$\Gamma$ -shaped domain (see figures 4 and 5) with vertices  $(0, 0)$ ,  $(0, 1)$ ,  $(1, 1)$ ,  $(-1, 1)$ ,  $(-1, -1)$  and  $(0, -1)$ .

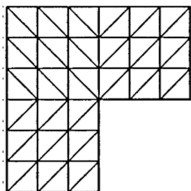


Figure:  $\Gamma$ -shape triangulation  
(without UBC)

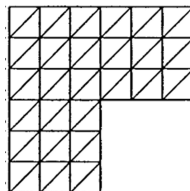


Figure:  $\Gamma$ -shape triangulation  
(with UBC)



## Model data:

### Finite element:

P1-Lagrange finite element.

### Meshsize:

The meshsize  $h_k$  for the  $k$ -th level grid is taken by  $2^{-k}$ .

### Multigrid parameters:

Using a W-cycle  $k$ -th ( $p=2$ ) level iteration, with 5 smoothing steps ( $m=n=5$ ). **Why  $m = n = 5$ ?** Because the numerical results do not appear to improve for any larger  $m$  or  $n$ .

Singular function on the  $\Gamma$ -shaped domain:

$$s_1(r, \theta) = \eta(r)r^{2/3} \sin(2/3\theta),$$

$$s_2(r, \theta) = \eta(r)r^{4/3} \sin(4/3\theta).$$

Cut-off function  $\eta$ :

$$\eta(r) = \begin{cases} 1 & 0 \leq r \leq \frac{1}{4} \\ -192r^5 + 480r^4 - 440r^3 + 180r^2 - \frac{135r}{4} + \frac{27}{8}, & \frac{1}{4} \leq r \leq \frac{3}{4} \\ 0 & \frac{3}{4} \leq r. \end{cases}$$

(see figure 6).

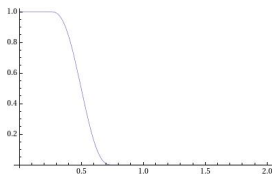


Figure: Cut off function  $\eta$

# Input data:

We will now compute a solution of the Poisson equation

$$\begin{aligned} -\Delta u &= f && \text{in } \Omega, \\ u &= 0 && \text{on } \partial\Omega. \end{aligned} \tag{13}$$

using

- Standard full multigrid algorithm.
- Full multigrid algorithm 1.
- Full multigrid algorithm 2.

for

$$f = -\Delta s_1 - \Delta s_2 + 6x(y^2 - y^4) + (x - x^3)(12y^2 - 2),$$

with exact solution

$$u = \underbrace{s_1}_{\in H^1(\Omega) \wedge \notin H^2(\Omega)} + \underbrace{s_2}_{\in H^2(\Omega) \wedge \notin H^{3-\epsilon}(\Omega)} + (x - x^3)(y^2 - y^4) \quad (\text{exact solution}).$$

# Experiment 1: Standard full multigrid algorithm (SFA)

Solving the Poisson equation (13)

$$-\Delta u = f \quad \text{in } \Omega, \quad u = 0 \quad \text{on } \partial\Omega.$$

with

$$f = -\Delta s_1 - \Delta s_2 + 6x(y^2 - y^4) + (x - x^3)(12y^2 - 2)$$

and exact solution

$$u = s_1 + s_2 + (x - x^3)(y^2 - y^4),$$

by the standard full multigrid algorithm, on the the  $\Gamma$ -shape, using the discretization fulfills **not the uniform band condition** presented in figure 4.

*Approximations for the stress intensity factors  $\kappa_k$  are computed by the extraction formula*

$$\kappa_h = \frac{1}{\pi} \left( \int_{\Omega} f s_- \, dx + \int_{\Omega} u \Delta s_- \, dx \right),$$

*using the P1 finite element solution  $u_k$  obtained by standard full multigrid algorithm.*

# Experiment 1: Standard full multigrid algorithm (SFA)

$e_k = |\Pi_k u - u_k|_{H^1(\Omega)}$  ... error in the energy norm,

$\sigma_k = \log_2 \left( \frac{|\kappa_{k-1} - 1|}{|\kappa_k - 1|} \right)$  ... convergence rate for stress intensity factor,

$\epsilon_k = \log_2 \left( \frac{e_{k-1}}{e_k} \right)$  ... convergence rate in the energy norm.

⇒ Theoretical:  $\sigma_k = \mathcal{O}(h_k^{4/3})$  and  $\epsilon_k = \mathcal{O}(h_k^{2/3})$

$k$	$\kappa_k$	$\sigma_k$	$e_k$	$\epsilon_k$
1	1.6999229601	–	$1.27093 \times 10^0$	–
2	1.2589102299	1.43	$5.91072 \times 10^{-1}$	1.1045
3	1.1036407706	1.32	$1.61387 \times 10^{-1}$	1.8728
4	1.0287080790	1.85	$5.74371 \times 10^{-2}$	1.4905
5	1.0073492045	1.97	$2.76732 \times 10^{-2}$	1.0535
6	1.0020544785	1.84	$1.64752 \times 10^{-2}$	0.7482
7	1.0005531037	1.89	$1.02811 \times 10^{-2}$	0.6803
8	1.0001571227	1.82	$6.46930 \times 10^{-2}$	0.6683
9	1.0000458701	1.78	$4.07502 \times 10^{-3}$	0.6668
10	1.0000142397	1.69	$2.56715 \times 10^{-3}$	0.6666
11	1.0000046460	1.62	$1.61722 \times 10^{-3}$	0.6666

Figure: Results for SFA.

## Experiment 2: Full multigrid algorithm 1 (FMGA1)

Solving the Poisson equation

$$-\Delta u = f \quad \text{in } \Omega, \quad u = 0 \quad \text{on } \partial\Omega.$$

with

$$f = -\Delta s_1 - \Delta s_2 + 6x(y^2 - y^4) + (x - x^3)(12y^2 - 2)$$

and exact solution

$$u = s_1 + s_2 + (x - x^3)(y^2 - y^4),$$

by the standard full multigrid algorithm 2, on the the  $\Gamma$ -shape, using the discretization **fulfills not the uniform band condition** presented in figure 4.

*We compute  $\kappa_k$  and  $w_k \in V_k$  which are approximations of ...*

- *stress intensity factor  $\kappa = 1$ ,*
- *and the regular part of the exact solution  $w = s_2 + (x - x^3)(y^2 - y^4)$ .*

# Experiment 2: Full multigrid algorithm 1 (FMGA1)

$e_k = |\Pi_k w - w_k|_{H^1(\Omega)}$  ... error in the energy norm,

$\sigma_k = \log_2 \left( \frac{|\kappa_{k-1} - 1|}{|\kappa_k - 1|} \right)$  ... convergence rate for stress intensity factor,

$\epsilon_k = \log_2 \left( \frac{e_{k-1}}{e_k} \right)$  ... convergence rate in the energy norm.

⇒ Theoretical:  $\sigma_k = \mathcal{O}(h_k^{5/3})$  and  $\epsilon_k = \mathcal{O}(h_k)$

$k$	$\kappa_k$	$\sigma_k$	$e_k$	$\epsilon_k$
1	-	-	$7.929 \times 10^{-1}$	-
2	1.69992296014	-	$8.364 \times 10^{-1}$	-0.07
3	0.82132136706	1.97	$2.322 \times 10^{-1}$	1.85
4	1.02037630458	3.13	$3.456 \times 10^{-2}$	2.75
5	0.99943755129	5.18	$6.236 \times 10^{-3}$	2.47
6	1.00003984026	3.82	$1.595 \times 10^{-3}$	1.97
7	1.00000536058	2.89	$4.200 \times 10^{-4}$	1.93
8	1.00000234005	1.20	$1.170 \times 10^{-4}$	1.84
9	1.00000057569	2.02	$3.567 \times 10^{-5}$	1.71
10	1.00000012632	2.19	$1.204 \times 10^{-5}$	1.57
11	1.00000002876	2.13	$4.397 \times 10^{-6}$	1.45
12	1.00000000746	1.95	-	-

Figure: Results for FMGA1.

## Experiment 3: Full multigrid algorithm 2 (FMGA2)

Solving the Poisson equation (13)

$$-\Delta u = f \quad \text{in } \Omega, \quad u = 0 \quad \text{on } \partial\Omega.$$

with

$$f = -\Delta s_1 - \Delta s_2 + 6x(y^2 - y^4) + (x - x^3)(12y^2 - 2)$$

and exact solution

$$u = s_1 + s_2 + (x - x^3)(y^2 - y^4),$$

by the standard full multigrid algorithm 2, on the the  $\Gamma$ -shape, using the discretization [fulfills the uniform band condition](#) presented in figure 5.

*We compute  $\kappa_{1,k}, \kappa_{2,k}$  and  $w_k \in V_k$  which are approximations of ...*

- *stress intensity factors  $\kappa_1 = \kappa_2 = 1$ ,*
- *and the regular part of the exact solution  $w = (x - x^3)(y^2 - y^4)$ .*



# Experiment 3: Full multigrid algorithm 2 (FMGA2)

$e_k = |\Pi_k w - w_k|_{H^1(\Omega)}$  ... error in the energy norm,

$\sigma_{i,k} = \log_2 \left( \frac{|\kappa_{i,k-1} - 1|}{|\kappa_{i,k} - 1|} \right)$  ... convergence rate for stress intensity factor,

$\epsilon_k = \log_2 \left( \frac{e_{k-1}}{e_k} \right)$  ... convergence rate in the energy norm.

⇒ Theoretical:  $\sigma_{i,k} = \mathcal{O}(h_k^2)$  and  $\epsilon_k = \mathcal{O}(h_k^2)$

$k$	$\kappa_{1,k}$	$\sigma_{1,k}$	$\kappa_{2,k}$	$\sigma_{2,k}$	$e_k$	$\epsilon_k$
1	–	–	–	–	$1.124 \times 10^0$	–
2	1.6229151283	–	1.17131298888	–	$7.361 \times 10^{-1}$	0.61
3	0.8859991798	2.45	0.99336080108	4.69	$1.417 \times 10^{-1}$	2.38
4	1.0091773397	3.63	1.00029662538	4.48	$1.131 \times 10^{-2}$	3.65
5	0.9999856171	9.32	1.00023130682	0.36	$5.829 \times 10^{-4}$	4.28
6	1.0000653041	–2.18	1.00002651087	3.13	$1.551 \times 10^{-4}$	1.91
7	1.0000136298	2.26	1.00000976600	1.44	$3.636 \times 10^{-5}$	2.09
8	1.0000044994	1.60	1.00000116447	3.07	$9.574 \times 10^{-6}$	1.93
9	1.0000011279	2.00	1.00000029598	1.98	$2.376 \times 10^{-6}$	2.01
10	1.0000002659	2.08	1.00000008791	1.75	$5.810 \times 10^{-7}$	2.03
11	1.0000000638	2.06	1.00000002475	1.82	$1.433 \times 10^{-7}$	2.02
12	1.0000000163	1.97	1.00000000585	2.08	–	–

Figure: Results for FMA2.

# Conclusions

- The multigrid methods use the simplest finite element.
- Since the grid are generated by connecting midpoints, it is easy to parallelize the algorithms.
- For more regular  $f$ , there exists a singular function representations where the regular part  $w$  is also more regular. In such cases multigrid methods with higher orders of convergence can be developed using higher order elements.
- Note that other superconvergence results which are less restrictive than the one based on the “uniform band” condition can also be used if they are available.