MULTI-GRID SOLUTIONS TO THE ELASTIC PLASTIC TORSION PROBLEM IN MULTIPLY CONNECTED DOMAINS

RONALD H. W. HOPPE

Technische Universität Berlin, FB Mathematik, Straße d, 17. Juni 135, D-1000 Berlin 12, West Germany

SUMMARY

The elastic plastic torsion problem for an elastic, perfectly plastic cylinder with multiply connected cross section twisted around its longitudinal axis is formulated as an obstacle problem for an associated stress potential, the obstacle being defined in terms of a generalized distance function. Based upon the reformulation of the obstacle problem as an equivalent linear complementarity problem, the latter is discretized by means of finite difference techniques, and a monotonically convergent iterative scheme for its numerical solution is developed. At each step of the iteration the solution of a reduced system of discrete Poisson equations is required which is done by applying multi-grid techniques with respect to a hierarchy of grid-point sets. Combined with a suitably chosen nested iteration process this results in a computationally very efficient algorithm for the approximate solution of the elastic plastic torsion problem.

1. INTRODUCTION

A widely used mathematical model for the problem to determine the equilibrium of an elastic, perfectly plastic medium is the Hencky model which allows the application of the powerful apparatus of convex analysis. However, it is well known that modelling elasto-plastic phenomena by Hencky's law has its limitations, the most severe one being that the history of work hardening of the material is not taken into account. For a more detailed discussion and for a general reference to the mathematical treatment of plasticity based on Hencky's model the reader is referred to Temam's textbook.

In particular, using the Hencky model, the elastic plastic torsion problem for an elastic, perfectly plastic cylinder with multiply connected cross section has been studied by Lanchon and Ting and numerically solved by Glowinski and Lanchon and Thomasset. The problem can be formulated as a constrained minimization problem for an associated energy functional or an equivalent variational inequality in terms of a related stress potential. Within this mathematical setup approximate solutions can be obtained using standard relaxation-projection techniques. However, it is well known that the convergence of these iterative methods crucially depends on a proper choice of the relaxation parameter. In other words, convergence may turn out to be rather slow if the relaxation parameter is not close to its optimal value. On the other hand, during the last decade 'optimal' iterative methods based on a recursive application of multiple grids have been developed for an efficient solution of algebraic systems arising from discretization of boundary-value problems for elliptic equations. In connection with variational inequalities such techniques have been successfully applied by Brandt and Cryer, Hackbusch and Mittelmann and by the author.
In this paper, we present a computational scheme for the approximate solution of the elastic plastic torsion problem based upon its reformulation as a linear complementarity problem which is given in Section 2. In Section 3 this problem is discretized by finite difference methods and a monotonically convergent iterative procedure for its solution is presented. Section 4 shows how the computational efficiency of multi-grid techniques can be taken into account both for the solution of the algebraic systems arising at each iteration step and for the generation of an appropriate start iterate. Finally, in Section 5 several numerical results are given for various geometrical configurations of the cylinder.

2. DEFINITION OF THE PROBLEM

We consider an elastic, perfectly plastic cylindrical bar \( Q := \Omega \times [0, l] \) of cross section \( \Omega \subset \mathbb{R}^2 \) and length \( l > 0 \) with \( m \) cylindrical cavities \( Q_i := \Omega_i \times [0, l] \) having the same direction of generatrices and cross section \( \Omega_i \subset \Omega \), \( 1 \leq i \leq m \), such that \( \Omega_1 \cap \Omega_j = \emptyset \), \( 1 \neq j \leq m \). Denoting by \( \partial Q_i := \Omega_i \times \{ l \} \), \( \partial Q_0 := \Omega \times \{ 0 \} \) the upper and lower ends respectively of the bar and by \( \partial Q_i := \Gamma \times (0, l) \), \( \Gamma = \partial \Omega \), the lateral surface, we suppose that at \( \partial Q_i \) the bar is twisted about the \( x_1 \)-axis by an angle \( \theta > 0 \) while \( \partial Q_0 \) is assumed to be stress-free.

For an isotropic material, using Hencky's law, modelling the plastic region according to the von Mises yield criterion and normalizing physical constants, it can be shown that the equilibrium stress tensor \( \{ \sigma_{ij} \} \) is of the form

\[
\sigma_{ij} = \begin{cases} 
\frac{\partial u_c}{\partial x_2} & \text{for } (i, j) = (1, 3) \text{ and } (i, j) = (3, 1) \\
-\frac{\partial u_c}{\partial x_1} & \text{for } (i, j) = (2, 3) \text{ and } (i, j) = (3, 2) \\
0 & \text{otherwise } (1 \leq i, j \leq 3)
\end{cases}
\]

The stress potential \( u_c \) is the unique solution of the constrained minimization problem

\[
J_C(u_c) = \inf_{v \in K} J_C(v)
\]

where \( J_C \) is the coercive, continuous, quadratic functional given by

\[
J_C(v) = \frac{1}{2} \int_{\Omega} |\nabla v|^2 \, dx - 2C \int_{\Omega} v \, dx
\]

\( C \) standing for the torsion angle per unit length, i.e. \( C = \theta / l \), and \( K \) is the closed, convex set

\[
K = \{ v \in H_0^1(\Omega) | v|_{\partial \Omega} = c_i = \text{const.}, 1 \leq i \leq m, |\nabla v| \leq 1 \text{ a.e. on } \Omega \}.
\]

Using standard arguments from convex analysis it follows that \( u_c \in K \) is the unique minimizer of \( J_C \) if and only if \( u_c \) satisfies the variational inequality

\[
u_c \in K
\]

\[
\int_{\Omega^*} \nabla u_c \cdot \nabla (v - u_c) \, dx \geq 2C \int_{\Omega^*} (v - u_c) \, dx, \quad v \in K
\]

where \( \Omega^* = \Omega \setminus \bigcup_{i=1}^{m} \bar{\Omega}_i \) is the region effectively occupied by the elastic plastic material. Note that
\( \Omega^* \) can be subdivided into the sets
\[
\Omega^*_{E,c} = \{ x \in \Omega^* \mid |\nabla u_c| < 1 \ \text{a.e.} \}
\]
\[
\Omega^*_{F,c} = \{ x \in \Omega^* \mid |\nabla u_c| = 1 \ \text{a.e.} \}
\]
characterizing the elastic and plastic region, respectively.

Setting \( \Omega_0 = \mathbb{R}^2 \setminus \Omega \) and \( \Gamma_0 = \partial \Omega_0 \) we define \( \Lambda_i, 0 \leq i \leq m, \) as the set of directed paths from \( \Gamma_i \) to \( \Gamma_0, \) a path \( P \in \Lambda_i \) consisting of directed edges \( P_{i_j,i_{j+1}}, 0 \leq j \leq n-1, n \in \mathbb{N}, \) of length \( \text{dist}((\Gamma_{i_j}, \Gamma_{i_{j+1}})) \) where \( i_0 = i \) and \( i_n = 0. \) Denoting by \( d_i \) the length of the shortest path within \( \Lambda_i, \) we define \( \psi : \Omega \to \mathbb{R} \) as the generalized distance function
\[
\psi(x) = \inf_{i \in \{0, 1, \ldots, m\}} \{ \text{dist}(x, \Omega_i) + d_i \}, \quad x \in \Omega
\]
Then, for \( C \to \infty \) it can be shown that the family \( \{u_c\} \) of solutions to (1) converges in \( H^1(\Omega) \) uniformly in \( C \) to \( u_\infty = \psi \) with \( |\nabla u_\infty| = 1 \ \text{a.e. in} \ \Omega^*. \) Consequently, \( \Omega^*_{E,\infty} = \emptyset, \ \Omega^*_{F,\infty} = \Omega^* \) and thus \( u_\infty \) represents the case of complete plastification.

The above fact motivates one to consider the modified minimization problem
\[
J_C(u_c) = \inf_{v \in K'} J_C(v)
\]
where
\[
K' = \{ v \in H^1_0(\Omega) \mid |v|_{\infty} = c_i, 1 \leq i \leq m, |v| \leq \psi \ \text{a.e. on} \ \Omega \}
\]
or the equivalent variational inequality
\[
 u_c \in K', \quad \int_{\Omega^*} \nabla u_c \cdot \nabla (v - u_c) \, dx \geq 2C \int_{\Omega^*} (v - u_c) \, dx, \quad v \in K'
\]
Indeed, for a simply connected domain \( \Omega, \) i.e. \( m = 0 \) and thus \( \psi(x) = \text{dist}(x, \Omega_0), x \in \Omega, \) it has been shown by Brézis and Sibony\(^3\) using a comparison principle for variational inequalities that problems (4a), (4b) and (10a), (10b) are equivalent. It is conjectured by Glowinski and Lanchon\(^5\) that this equivalence also holds true in the multiply connected case, a conjecture which is supported by various numerical results obtained via the approximate solution of (1) and (8), respectively.

From now on we will assume \( C > 0. \) In this case we have \( u_c > 0 \ \text{a.e. on} \ \Omega \) and hence, the constraint \( |u_c| \leq \psi \ \text{a.e. on} \ \Omega \) in the definition of \( K' \) can be replaced by \( 0 \leq u_c \leq \psi \ \text{a.e. on} \ \Omega. \) Then, taking advantage of the regularity result\(^1\)
\[
 u_c \in H^1_0(\Omega) \cap C^\infty(\Omega^*) \cap C^{1,\eta}(\overline{\Omega}^*), \quad x \in (0, 1)
\]
which holds true under the assumption of piecewise smooth \( \Gamma_i, 0 \leq i \leq m, \) and using the fact that the lower obstacle is never active, it can be easily shown that (10a), (10b) is equivalent to the linear complementarity (LC-) problem
\[
-(\Delta u_c)(x) \leq 2C, \quad u_c(x) \leq \psi(x), \quad x \in \Omega^*
\]
\[
(\Delta u_c)(x) + 2C(u_c(x) - \psi(x)) = 0, \quad x \in \Omega^*
\]
\[
u_c(x) = 0, \quad x \in \Gamma, \quad u_c(x) = c_i, \quad x \in \Gamma_i, \quad 1 \leq i \leq m
\]
The above LC-problem will serve as the basis for a computational scheme which will be developed in the following section.

3. THE ITERATIVE SCHEME

For the approximate solution of the LC-problem (12) we choose standard finite difference techniques with respect to a two-dimensional grid which for notational convenience will be assumed to be equidistant.

For a given step-size $h > 0$ we set

$$\mathbb{R}_h^2 = \{ z = (z_1, v_1, z_2, v_2) \mid z_i, v_i \in \mathbb{R}, \quad i = 1, 2 \}$$

and we define

$$\Omega_h = \mathbb{R}_h^2 \cap \Omega, \quad \Omega_{h,i} = \mathbb{R}_h^2 \cap \Omega_i, \quad 1 \leq i \leq m$$

$$\Omega_{h}^{*} = \mathbb{R}_h^{*} \setminus \bigcup_{i=1}^{m} \Omega_{h,i}$$

$\Omega_{h}^{*}$ will be called the set of interior grid-points.

Setting $e_1 = (1, 0), e_2 = (0, 1), e_3 = e_1 + e_2$ and $e_4 = e_2 - e_1$, with each $z = (z_1, z_2) \in \Omega_{h}^{*}$ we associate eight neighbouring grid-points which are situated in the directions $e_1, e_2, e_3, e_4, e_5, e_6, e_7, e_8$, respectively, and the associated grid-points will be constructed in the following way.

Let us denote by $L_{SW}(z)$ the line segment

$$L_{SW}(z) = \{ x = (x_1, x_2) \in \mathbb{R}^2 \mid x_2 = h^{-1}x_1 + (z_2 - h^{-1}z_1), \quad z_1 - h \leq x_1 \leq z_1 \}$$

Then, if $L_{SW}(z) \subset \Omega^{*}$, we set $z_{SW} = (z_{SW,1}, z_{SW,2}), z_{SW,\mu} = z_{\mu} - h, 1 \leq \mu \leq 2$. Otherwise, we consider the non-empty set $\Gamma^{*} \cap L_{SW}(z)$, and we define $z_{SW}$ as the point in $\Gamma^{*} \cap L_{SW}(z)$ which has the shortest distance to $z$, i.e. $\text{dist}(z, z_{SW}) = \text{dist}(z, \Gamma^{*} \cap L_{SW}(z))$.

The line segments $L_S(z), L_{SE}(z), \ldots, L_{NE}(z)$ and the corresponding neighbours $z_S, z_{SE}, \ldots, z_{NE}$ will be constructed analogously.

We remark that for discretization with respect to a fixed grid only the four neighbouring points $z_S, z_W, z_E$ and $z_N$ corresponding to the standard five-point discretization of the Laplacian will be important while the neighbours $z_{SW}, z_{SE}, z_{NW}$ and $z_{NE}$ come into effect in the nested iteration process of a multi-grid procedure which will be described in detail in the next section.

Accordingly, we define the following sets of neighbouring grid-points

$$N_h^1(z) = \{ z_S, z_W, z_E, z_N \}$$

$$N_h^2(z) = N_h^1(z) \cup \{ z_{SW}, z_{SE}, z_{NW}, z_{NE} \}$$

and the corresponding grid-point sets

$$\Gamma_h^{*,\mu} = \left( \bigcup_{z \in \Omega_h^{*}} N_h^\mu(z) \right) \cap \Gamma^{*}, \quad 1 \leq \mu \leq 2$$

$$\Omega_h^{*,\mu} = \Omega_h^{*} \cup \Gamma_h^{*,\mu}, \quad 1 \leq \mu \leq 2$$

Specifically, we set

$$\Gamma_h^{*,\mu} = \Gamma_h^{*,\mu} \cap \Gamma_i, \quad 0 \leq i \leq m, \quad 1 \leq \mu \leq 2$$
The elements of $\Gamma_h^{*,\nu}$ and $\Gamma_h^{*,1,\nu}$ are called boundary grid-points while an interior grid-point $z \in \Omega_h^*$ is said to be regular if $N^h_1(z) \subset \Omega_h^*$, and irregular, otherwise.

We define $C(\Omega_h^*)$, $C(\Omega_h^{*,\mu})$ and $C(\Omega_h^{*,1,\mu})$, $1 \leq \mu \leq 2$, as the vector space of real-valued grid-functions on $\Omega_h^*$, $\Omega_h^{*,\mu}$ and $\Omega_h^{*,1,\mu}$ respectively. Moreover, $C_b(\Omega_h^{*,\mu})$, $1 \leq \mu \leq 2$, refers to the subspace of $C(\Omega_h^{*,\mu})$ given by

$$C_b(\Omega_h^{*,\mu}) = \{ u_h \in C(\Omega_h^{*,\mu}) \mid u_h(z) = 0, \quad z \in \Gamma_h^{*,\mu}, \quad 1 \leq i \leq m \}$$

For a grid-function $u_h \in C(\Omega_h^{*,1})$ and $z \in \Omega_h^*$ we define the standard forward and backward difference quotients respectively

\begin{align*}
(D_{h^+} u_h)(z) &= \frac{u_h(z) - u_h(z_j)}{z_j - z_i}, \quad 1 \leq j \leq 2 \\
(D_{h^-} u_h)(z) &= \frac{u_h(z) - u_h(z_j)}{z_j - z_i}, \quad 1 \leq j \leq 2
\end{align*}

where $z = z_j$ for $j = 1$, $z = z_N$ for $j = 2$ in (20a) and $z = z_{j-1}$ for $j = 1$, $z = z_2$ for $j = 2$ in (20b).

Further we define

$$\delta^2 u_h(z) = (D_{h^+} u_h)(z) + (D_{h^-} u_h)(z) \quad \text{as an approximation of the Laplacian.}$$

Note that for a regular interior grid-point (21) reduces to the standard five-point discretization with respect to an equidistant grid while for an irregular interior grid-point (21) corresponds to the well-known Shortley–Weller scheme.

Introducing a discrete obstacle $\psi_h \in C(\Omega_h^*)$ by $\psi_h = \psi|_{\Omega_h^*}$, we consider the discrete LC-problem

\begin{align*}
-(\delta^2 u_{h,c})(z) &\leq 2C, \quad u_{h,c}(z) \leq \psi_h(z), \quad z \in \Omega_h^* \quad \text{(22a)} \\
\left(\delta^2 u_{h,c}(z) + 2C\right)(u_{h,c}(z) - \psi_h(z)) &= 0, \quad z \in \Omega_h^* \quad \text{(22b)} \\
u_{h,c}(z) &= 0, \quad z \in \Gamma_h^{*,1}, \quad u_{h,c}(z) = c_i, \quad z \in \Gamma_h^{*,0}, 1 \leq i \leq m \quad \text{(22c)}
\end{align*}

We can formally rewrite (22) as the Hamilton–Jacobi–Bellman (HJB-) equation

\begin{align*}
\max \left[ -(\delta^2 u_{h,c})(z) - 2C, \quad u_{h,c}(z) - \psi_h(z) \right] &= 0, \quad z \in \Omega_h^* \quad \text{(23a)} \\
u_{h,c}(z) &= 0, \quad z \in \Gamma_h^{*,1}, \quad u_{h,c}(z) = c_i, \quad z \in \Gamma_h^{*,0}, 1 \leq i \leq m \quad \text{(23b)}
\end{align*}

Assuming that for increasing $C$ the plastification spreads from $\Gamma$, $0 \leq i \leq m$, it is natural to take $c_i = d_i$, $1 \leq i \leq m$. Then, setting $N_h = \text{card } \Omega_h^*$, (23) is a system of $N_h$ non-linear difference equations in the $N_h$ unknowns $u_h(z), z \in \Omega_h^*$. Following the arguments used for discrete LC-problems and discrete HJB-equations\textsuperscript{10,13} it can be easily shown that (22) or (23) admit a unique solution $u_{h,c} \in C_b(\Omega_h^{*,1})$ which approximates the solution $u_c$ to (12) of order $O(h^2)$ provided $u_c \in C^2(\Omega^*)$, $x \in (0, 1)$. Moreover, in view of the regularity result (11) interior estimates of order $O(h^2)$ can be expected.

We now present a monotonic convergent iterative scheme for the solution of the discrete LC-problem (22) which has been successfully used for discrete HJB-equations.\textsuperscript{10,13} At each iteration step this scheme requires the solution of a reduced linear algebraic system which can be constructed in the following way.

Given an iterate $u_{h,c}^* \in C_b(\Omega_h^{*,1})$, $v \geq 0$, we set

$$d^{(v)}_{h,c}(z) = -(\delta^2 u_{h,c}^*)(z) - 2C, \quad d^{(v)}_{h,c}(z) = u_{h,c}^*(z) - \psi_h(z),$$

$z \in \Omega_h^*$, and we define a difference operator $A^v_{h,c} : C_b(\Omega_h^{*,1}) \to C(\Omega_h^*)$ and a grid-function $f^{(v)}_{h,c} \in C(\Omega_h^*)$...
by

\[
(A_h^{(v)} v_h)(z) = \begin{cases} 
(-\delta_h v_h)(z), & \text{if } d_h^{(v)}(z) > d_h^{(v)}(z) \\
v_h(z), & \text{otherwise}
\end{cases}
\]

(24)

\[
f_h^{(v)}(z) = \begin{cases} 
2C, & \text{if } d_h^{(v)}(z) > d_h^{(v)}(z) \\
\psi_h(z), & \text{otherwise}
\end{cases}
\]

(25)

Then, a new iterate \( u_h^{*+1} \in C(u(\Omega_h^{*+1})) \) will be computed as the solution of

\[
(A_h^{(v)} u_h^{*+1})(z) = f_h^{(v)}(z), \quad z \in \Omega_h^{*}
\]

(26)

The definitions of \( A_h^{(v)} \) and \( f_h^{(v)} \) induce a partition of \( \Omega_h^{*} \) according to

\[
\Omega_h^{*} = \Omega_h^{*(1)}(u_h^{*+1}, C) \cup \Omega_h^{*(2)}(u_h^{*+1}, C), \quad \Omega_h^{*(1)}(u_h^{*+1}, C) \cap \Omega_h^{*(2)}(u_h^{*+1}, C) = \emptyset
\]

(27)

where

\[
\Omega_h^{*(1)}(u_h^{*+1}) = \{ z \in \Omega_h^{*} | d_h^{(1), (v)}(z) > d_h^{(2), (v)}(z) \}
\]

(28a)

\[
\Omega_h^{*(2)}(u_h^{*+1}) = \{ z \in \Omega_h^{*} | d_h^{(1), (v)}(z) \leq d_h^{(2), (v)}(z) \}
\]

(28b)

Setting \( N_h = \text{card} \Omega_h^{*(u)}(u_h^{*+1}), 1 \leq \mu \leq 2 \), we have \( u_h^{*+1} = \psi_h(z), z \in \Omega_h^{*(2)}(u_h^{*+1}, C) \), and hence (26) reduces to a system of \( N_h \) linear equations in the \( N_h \) unknowns \( u_h^{*+1}(z), z \in \Omega_h^{*} (u_h^{*+1}) \).

Since \( \Omega_h^{*(2)}(u_h^{*+1}) \) represents the set of grid-points where the discrete obstacle function \( \psi_h \) is active, grid-points \( z \in \Omega_h^{*(2)}(u_h^{*+1}, C) \) are said to be active while grid-points \( z \in \Omega_h^{*(1)}(u_h^{*+1}, C) \) are called inactive. For the exact solution \( u_h, C \) to (22) or (23) the set \( \Omega_h^{*(2)}(u_h, C) \) characterizes the discrete coincidence set which indicates the discrete region of plastification.

Denoting by \( \| \cdot \|_{\infty} \) the maximum norm in \( C(\Omega_h^{*}) \) and prespecifying an accuracy bound \( \varepsilon_h^1 > 0 \) as termination criterion for the iteration, the complete algorithm \( \text{EPT}(u_h^{*+1}, C) \) for the solution of the elastic plastic torsion problem with the torsion angle per unit length \( C \) and the accuracy bound \( \varepsilon_h^1 \) as inputs and an approximate solution \( u_h^{*+1} \) as output reads as follows.

ALGORITHM \( \text{EPT}(u_h^{*+1}, \varepsilon_h^1, C) \):

Step 1. Choose a start-iterate \( u_h^0 \in C_u(\Omega_h^{*+1}) \) and set \( v = 0 \).

Step 2. Determine \( A_h^{(v)} \) and \( f_h^{(v)} \) according to (24), (25) and compute \( u_h^{*+1} \in C_u(\Omega_h^{*+1}) \) as the solution to (26).

Step 3. If \( \| u_h^{*+1} - u_h^{*+1} \|_{\infty} < \varepsilon_h^1 \), then set \( u_h, C = u_h^{*+1} \) and stop the algorithm, otherwise set \( v = v + 1 \) and go to Step 2.

Recalling that a grid-function \( w_h \in C_u(\Omega_h^{*+1}) \) is said to be a supersolution to (23) if

\[
\max \left[ (-\delta_h w_h)(z) - 2C, w_h(z) - \psi_h(z) \right] \geq 0, \quad z \in \Omega_h^{*}
\]

(29)

we have the following convergence result.\(^{11}\)

**Theorem**

For given \( C > 0 \) let \( u_h^0 \in C_u(\Omega_h^{*+1}) \) be a supersolution to (23). Then, the sequence \( \{ u_h^{*+1} \}_{v \geq 0} \) of iterates \( u_h^{*+1} \in C_u(\Omega_h^{*+1}) \) governed by the algorithm \( \text{EPT}(u_h, C, \varepsilon_h^1) \) is a monotonically decreasing sequence of grid-functions converging for \( v \to \infty \) to the unique solution \( u_h, C \) of (23).

**Proof.** For \( v \geq 1 \) we have

\[
\max \left[ (-\delta_h u_h^{*+1})(z) - 2C, u_h^{*+1}(z) - \psi_h(z) \right] \geq (A_h^{(v-1)} u_h^{*+1})(z) - f_h^{(v-1)}(z) = 0, \quad z \in \Omega_h^{*}
\]

(30)
and hence, each iterate \( u_h^{v,c}, v \geq 0 \), is a supersolution to (23). Consequently,

\[
(A_h^{(v-1)} u_h^{(v-1)}(z) - f_h^{(v-1)}(z)) \geq 0, \quad z \in \Omega_h^*
\]

Moreover,

\[
0 = \max \left[ (-\delta_h^2 u_h^{v,c}(z) - 2C, u_h^{v,c}(z) - \psi_h(z) \right]
\geq (A_h^{(v-1)} u_h^{v,c}(z) - f_h^{(v-1)}(z), \quad z \in \Omega_h^*
\]

Since for each \( v \geq 1 \) the operator \( A_h^{(v-1)} \) satisfies the maximum principle

\[
A_h^{(v-1)} w_h = A_h^{(v-1)} w_h = v_h \geq w_h, \quad v_h, w_h \in C_0(\Omega_h^{v-1})
\]

inequalities (30), (31) and (32) yield \( u_h^{v,c} \geq u_h^{v,c} \geq u_h^{v,c}, v \geq 1 \). This implies the existence of a grid-function \( u_h^{v,c} \in C_0(\Omega_h^{v-1}) \) such that \( u_h^{v,c}(z) \rightarrow u_h^{v,c}(z), z \in \Omega_h^* \), for \( v \rightarrow \infty \). In view of

\[
\max \left[ (-\delta_h^2 u_h^{v-1,c}(z) - 2C, u_h^{v-1,c}(z) - \psi_h(z) \right]
\geq \left(A_h^{(v-1)} u_h^{v-1,c}(z) - f_h^{(v-1)}(z) = (A_h^{(v-1)}(u_h^{v-1,c} - u_h^{v,c}))(z)\right)
\]

for \( v \rightarrow \infty \) we get

\[
\max \left[ (-\delta_h^2 u_h^{v,c}(z) - 2C, u_h^{v,c}(z) - \psi_h(z) \right] = 0, \quad z \in \Omega_h^*
\]

whence \( u_h^{v,c} = u_h^{v,c} \) because of the unique solvability of (23).

Bearing in mind that the solution \( u_h^{v,c} \) to (23) satisfies the strict complementary slackness condition

\[
(-\delta_h^2 u_h^{v,c}(z) - 2C, z \in \Omega_h^{v,2}(u_h^{v,c})
\]

it can be shown that convergence occurs after a finite number of steps. In particular, we have the following.

**Corollary**

Under the assumptions of the preceding theorem there exist indices \( v_0, v_1 \in \mathbb{N}, v_1 \geq v_0 \), such that

\[
\Omega_h^{v,0}(u_h^{v,c}) \subseteq \Omega_h^{v,1}(u_h^{v,c}) \subseteq \Omega_h^{v,2}(u_h^{v,c}), \quad v \geq v_0
\]

and

\[
\Omega_h^{v,2}(u_h^{v,c}) = \Omega_h^{v,2}(u_h^{v,c}), \quad u_h^{v,c} = u_h^{v,c}, \quad v \geq v_1
\]

**Proof.** The proof is by contradiction arguments. If the first inclusion in (34) does not hold, there exist a subsequence \( \mathbb{N}' \subset \mathbb{N} \) and a grid-point \( z \in \Omega_h^* \) with \( z \notin \Omega_h^{v,2}(u_h^{v,c}) \) but \( z \notin \Omega_h^{v,2}(u_h^{v,c}) \), \( v \in \mathbb{N}' \). Consequently, \( (-\delta_h^2 u_h^{v,c}(z) - 2C < 0 \) and \( (-\delta_h^2 u_h^{v,c}(z) - 2C = 0, v \in \mathbb{N}', \) contradicting the fact that \( (-\delta_h^2 u_h^{v,c}(z) - 2C \rightarrow (-\delta_h^2 u_h^{v,c}(z) - 2C \rightarrow \infty, v \in \mathbb{N}' \).

Similarly, if the second inclusion is not true, there exist another subsequence \( \mathbb{N}'' \subset \mathbb{N} \) and a grid-point \( z \in \Omega_h^* \) such that \( z \notin \Omega_h^{v,2}(u_h^{v,c}) \) but \( z \notin \Omega_h^{v,2}(u_h^{v,c}) \), \( v \in \mathbb{N}'' \), whence

\[
(-\delta_h^2 u_h^{v,c}(z) - 2C > u_h^{v,c}(z) - \psi_h(z) \geq u_h^{v,c}(z) - \psi_h(z) \geq u_h^{v,c}(z) - \psi_h(z) \geq (-\delta_h^2 u_h^{v,c}(z) - 2C = 0
\]

For \( v \rightarrow \infty \) this gives \( u_h^{v,c}(z) - \psi_h(z) = 0 \) and \( (-\delta_h^2 u_h^{v,c}(z) - 2C = 0, \) contradicting the strict complementary slackness condition (33).

Finally, since there is only a finite number of constraints, there must be an integer \( v_1 \geq v_0 \) such that (34) holds true and thus \( u_h^{v,c} = u_h^{v,c} \).
The assumption on the start-iterate \( u_{h,c}^0 \in C_b(\Omega^{*}_{k+1}) \) to be a supersolution to (23) can be dispensed with. Indeed, if \( u_{h,c}^0 \in C_b(\Omega^{*}_{k+1}) \) is chosen arbitrarily, then it is guaranteed that after one iteration we end up with a supersolution \( u_{h,c}^1 \), and thus we obtain monotonic convergence of the sequence \( \{u_{h,c}^i\}_{i \geq 1} \). When using multiple grids this global convergence property is of importance in the so-called nested iteration process, as will be described in the next section.

### 4. THE MULTI-GRID ALGORITHMS

As has been shown in the last section, the implementation of the iterative procedure \( \text{EPT}(u_{h,c}, \mathcal{E}_h, C) \) requires the successive solution of discrete Poisson equations with respect to an increasing sequence of grid-point sets \( \Omega^{*}_{k+1}(u_{h,c}^i), \forall i \in \mathbb{N} \). Since in view of the geometry of these sets fast-Poisson solvers are not applicable, and since standard iterative solvers like Gauss-Seidel overrelaxation suffer from poor convergence, we recommend the application of multi-grid techniques with respect to a hierarchy \( \{\Omega^{*}_{k}\}_{k=0}^{l-1} \) of equidistant grids with step-sizes \( h_{k+1} < h_k, 0 \leq k \leq l-1 \). For simplicity we assume \( h_{k+1} = h_k/2, 0 \leq k \leq l-1 \), given some \( h_0 > 0 \).

By \( \Omega^{*}_{k+1}(u_{h,c}^i) \) and \( \Omega^{*}_{k+1}(u_{h,c}^i) \) (1 \( \leq k \leq l \)), we denote the grids defined according to (14), (17) while \( \delta_{k+1}^{(2)}(u_{h,c}^i), 0 \leq k \leq l \), refers to the discrete Laplacian on \( C_b(\Omega^{*}_{k+1}) \) constructed by means of (21). Then, in contrast to standard multi-grid methods for discrete Poisson equations, the problem which arises in connection with equation (26) on the finest grid \( \Omega^{*}_{1} \) is how to specify the sets \( \Omega^{*}_{k+1}(u_{h,c}^i), 1 \leq k \leq l \), of active/inactive grid-points on the lower levels \( 0 \leq k < l \). In order to avoid that discrete free boundaries remain undetected it is recommended to specify an interior grid-point \( z \in \Omega^{*}_{k}, k < l, \) as inactive, if \( z \) and all its eight neighbours from \( N^2(z) \) are inactive on level \( k+1. \) This means, given \( \Omega^{*}_{k+1}(u_{h,c}^i), 0 \leq k \leq l \), we define recursively

\[
\Omega^{*}_{k+1}(u_{h,c}^i) = \Omega^{*}_{k+1}(u_{h,c}^i) \cap \{z \in \Omega^{*}_{k+1}(u_{h,c}^i) \cap N^2(z) < \Omega^{*}_{k+1}(u_{h,c}^i) \}
\]

We set \( \Omega^{*}_{k+1}(u_{h,c}^i) = \Omega^{*}_{k+1}(u_{h,c}^i) \cap \Omega^{*}_{k+1}(u_{h,c}^i), 0 \leq k \leq l \), and we define \( k_{\min} \) as the smallest integer \( 0 \leq k \leq l \) for which \( \Omega^{*}_{k+1}(u_{h,c}^i) \neq \emptyset \), i.e. \( \Omega^{*}_{k_{\min}} \) is the coarsest grid with at least one inactive grid-point.

For the coarse-to-fine and the fine-to-coarse transfers within a multi-grid cycle we have to provide suitable prolongations and restrictions. We denote by \( p^k_{k+1} : C_b(\Omega^{*}_{k}) \to C_b(\Omega^{*}_{k+1}), 0 \leq k \leq l-1 \), the prolongation based upon bilinear interpolation and by \( r^k_{k+1} : C_b(\Omega^{*}_{k+1}) \to C_b(\Omega^{*}_{k}), 0 \leq k \leq l-1 \), the corresponding full weighted restriction. Further, if \( v_k \in C_b(\Omega^{*}_{k}) \) or \( v_k \in C_b(\Omega^{*}_{k+1}), 0 \leq k < l \), we introduce a grid-function \( \tilde{v}_k \in C_b(\Omega^{*}_{k}) \) by

\[
\tilde{v}_k(z) = \begin{cases} v_k(z), & \text{if } z \in \Omega^{*}_{k+1}(u_{h,c}^i) \\ 0, & \text{otherwise} \end{cases}
\]

Then, we define prolongations \( p^k_{k+1} : C_b(\Omega^{*}_{k+1}) \to C(\Omega^{*}_{k+1}) \) and restrictions \( r^k_{k+1} : C(\Omega^{*}_{k+1}) \to C(\Omega^{*}_{k}) \) or \( p^k_{k+1} : C_b(\Omega^{*}_{k+1}) \to C(\Omega^{*}_{k+1}) \) and \( r^k_{k+1} : C_b(\Omega^{*}_{k+1}) \to C(\Omega^{*}_{k+1}) \), \( 0 \leq k \leq l-1 \), according to

\[
(p^k_{k+1} v_k)(z) = \begin{cases} (p^k_{k+1} \tilde{v}_k)(z), & \text{if } z \in \Omega^{*}_{k+1}(u_{h,c}^i) \\ 0, & \text{otherwise} \end{cases}
\]

\[
[r^k_{k+1} v_{k+1}](z) = \begin{cases} (r^k_{k+1} \tilde{v}_{k+1})(z), & \text{if } z \in \Omega^{*}_{k+1}(u_{h,c}^i) \\ 0, & \text{otherwise} \end{cases}
\]

Further, we need a suitable smoothing procedure for the reduced linear system (26) on level \( k = l \) and for the defect correction equations on levels \( k_{\min} < k < l \) as well as a solver for the defect correction equation on the coarsest grid. For both purposes we propose Gauss–Seidel iteration, denoting the application of \( k \) Gauss–Seidel iterations to the corresponding equation on level \( k \) by

\[
w_k = S_k(v_k, k)
\]
MULTIPLY CONNECTED DOMAINS

where \( v_k \in C_b(\Omega^*_{k+1}) \) is used as a start-iterate. Note that the defect correction equations on levels \( k_{\text{min}} \leq k < l \) are given by

\[
(A^{(i)}_k u_{k+1}^{(i)})_k(z) - f^{(i)}_k(z) = 0, \quad z \in \Omega^*_k
\]

where the operator \( A^{(i)}_k : C_b(\Omega^*_{k+1}) \to C(\Omega^*_k) \) and the grid-function \( f^{(i)}_k \in C(\Omega^*_k) \) are defined according to

\[
(A^{(i)}_k v_k)(z) = \begin{cases} (-\bar{\delta}_k^2 v_k)_k(z), & \text{if } z \in \Omega^*_{k}^{(1)}(u_0^c) \\ v_k(z), & \text{if } z \in \Omega^*_{k}^{(2)}(u_0^c) \end{cases}
\]

\[
f^{(i)}_k(z) = (A^{(i)}_k v_k(z)) - (\bar{\delta}_k^2 (A^{(i)}_{k+1} \bar{u}_{k+1}^c - f^{(i)}_{k+1}))(z)
\]

\( \bar{u}_{k+1}^c \) denoting the smoothed iterate on level \( k+1 \) and \( v_k^c \in C_b(\Omega^*_{k+1}) \) given by

\[
v_k^c(z) = \begin{cases} (\bar{\delta}_k^2 \bar{u}_{k+1}^c)_k(z), & \text{if } z \in \Omega^*_{k}^{(1)}(u_0^c) \\ \psi_k(z), & \text{if } z \in \Omega^*_{k}^{(2)}(u_0^c) \\ d_i, & \text{if } z \in \Gamma^c_{k,i}(1 \leq i \leq m) \\ 0, & \text{if } z \in \Gamma^c_{k,0} \end{cases}
\]

If \( k = l \), we set \( v_k^c = u_k^c \).

The following algorithm \( \text{MG}(u_{k+1}^c, v_k^c, k) \) describes a complete multi-grid cycle for computation of an approximation \( u_{k+1}^c \in C_b(\Omega^*_{k+1}) \) to (26), if \( k = l \), or to (40), if \( k < l \), given \( v_k^c \in C_b(\Omega^*_{k+1}) \) as a start-iterate.

ALGORITHM \( \text{MG}(u_{k+1}^c, v_k^c, k) \):

Step 1. If \( k = k_{\text{min}} \), compute \( u_{k+1}^c \) by \( \kappa_3 \) Gauss–Seidel iterations applied to (40) on level \( k_{\text{min}} \), i.e.

\[
u_{k+1}^c = S_k(v_k^c; \kappa_3)
\]

Step 2. If \( k > k_{\text{min}} \), compute a smoothed iterate \( \bar{u}_k^c \) by \( \kappa_1 \) Gauss–Seidel iterations applied to (26), if \( k = l \), or to (40), if \( k < l \), i.e.

\[
\bar{u}_k^c = S_k(v_k^c; \kappa_1)
\]

Step 3. Determine the operator \( A^{(i)}_{k-1} \) and the grid-functions \( f^{(i)}_{k-1} \), \( v_{k-1}^c \) by means of (41), (42), (43) and compute \( u_{k-1}^c \) by applying \( \gamma_k \) cycles \( \text{MG}(u_{k-1}^c, v_{k-1}^c, k-1) \).

Step 4. Compute \( \bar{u}_{k-1}^c \) by \( C_b(\Omega^*_{k+1}) \) by

\[
\bar{u}_{k-1}^c = \bar{u}_{k-1}^c - \bar{\delta}_{k-1}^2 [\bar{\delta}_k^{k-1} \bar{u}_{k-1}^c - u_{k-1}^c] + \bar{\delta}_{k-1}^2 [\bar{\delta}_k^{k-1} \bar{u}_{k-1}^c - u_{k-1}^c]
\]

and determine \( u_{k-1}^c \) by \( \kappa_2 \) Gauss–Seidel iterations applied to (26), if \( k = l \), or to (40), if \( k < l \), using \( \bar{u}_{k-1}^c \) as start-iterate, i.e.

\[
u_{k-1}^c = S_k(\bar{u}_{k-1}^c; \kappa_2)
\]

Finally, if \( k < l \), set \( v_k^c = u_{k+1}^c \).

Due to the definition of \( \bar{\delta}_k^{k+1} \) and \( \bar{\delta}_k^{k+1} \), and the fact that the smoothing procedure only changes values in \( z \in \Omega^*_{k}^{(1)}(u_0^c) \), it follows from (45), (46) and (47) that the prespecified values \( \psi_k(z) \) in \( z \in \Omega^*_{k}^{(2)}(u_0^c) \) remain unchanged while performing a multi-grid cycle.

Choosing an accuracy bound \( \epsilon_t > 0 \) as termination criterion for the multi-grid iterations, the iterative procedure EPT of Section 3 can be coupled with the multi-grid scheme MG to give an effective algorithm \( \text{MG-EPT}(u_0^c, \epsilon_t^0, \epsilon_t^c, C) \) for computing an approximation \( u_l^c \in C_b(\Omega^*_{l+1}) \) to the LC-problem (12) on level \( l \). This algorithm results if in EPT(\( u_k^c, \epsilon_l^0, \epsilon_l^c, C \)) we take \( h = h_l \) and replace Step 2 by
Step 2'. (i) Determine $A^{(n)}, f^{(n)}$ and the sets $\Omega_k^{\bullet, (n)}(u^0_{l,c})$, $1 \leq \eta \leq 2$, $k_{\min} \leq k \leq l$, according to (24), (25), (28), (36), set $u^{l+1,0}_{o,c} = u^0_{l,c}$ and $\mu = 0$.

(ii) Set $\mu = \mu + 1$ and determine $u^{l+1,\mu}_{l,c}$ by performing $\text{MG}(u^{l+1,\mu}_{l,c}, v^l_{o,c}, l)$.

(iii) If $\|u^{l+1,\mu}_{l,c} - u^{l+1,\mu-1}_{l,c}\|_{\infty} < \varepsilon^l$, set $u^{l+1,\mu}_{l,c} = u^{l+1,\mu-1}_{l,c}$ and go to Step 3, otherwise set $u^{l+1,\mu}_{l,c} = u^{l+1,\mu}_{l,c}$ and go to Step 2 (ii).

The performance of the multi-grid algorithm MGEPT can be considerably improved by choosing an appropriate start-iterate $u^0_{l,c}$ on the finest grid. This can be achieved by nested iteration. 

Starting from a grid-function $u^0_{0,c}$ on the coarsest grid, at each level $0 \leq k < l$ an approximation $u^*_k$ is computed by performing a multi-grid cycle involving the grids $0 < m < k$, and $u^*_k$ is then interpolated to the level $k + 1$ by a suitable prolongation process. For this purpose we cannot use the prolongation operator $\beta^k$ as defined by (37), since $\beta^k$ is specially designed for the coarse-to-fine transfer within MG already assuming the knowledge of an approximation on level $k + 1$. Therefore, we construct another prolongation operator $\beta^k$:

If $u^k \in C_0(\Omega^*_k)$ and $z \in \Omega^*_k \setminus \Omega^*_k$, we define $\delta_k(z')$, $z' \in N^2_k(z)$, by

$$
\delta_k(z') = \begin{cases} 
    v_k(z'), & \text{if } z' \in \Omega^*_k \\
    0, & \text{if } z' \in \Gamma^*_k+1,0 \\
    d_i, & \text{if } z' \in \Gamma^*_k+1,i, \quad 1 \leq i \leq m
\end{cases}
$$

where (i) $z' = z_{SW}$, $z_{SE}$, if $z = h_{k+1} e_{1} \in \mathbb{N}^2_k$, (ii) $z' = z_{SN}$, $z_{SW}$, if $z = h_{k+1} e_{2} \in \mathbb{N}^2_k$, and (iii) $z' = z_{SW}$, $z_{SE}$, $z_{NW}$, $z_{NE}$, if $z = h_{k+1} e_{3} \in \mathbb{N}^2_k$, $\mu = 3, 4$.

We then define $(\beta^k v_k)(z)$ by $(\beta^k v_k)(z) = v_k(z)$, if $z \in \Omega^*_k$, and, if $z \notin \Omega^*_k$, by linear interpolation between $\delta_k(z_{SW})$ and $\delta_k(z_{E})$ in case (i) or $\delta_k(z_{S})$, $\delta_k(z_{N})$ in case (ii), and by bilinear interpolation between $\delta_k(z_{SW})$, $\delta_k(z_{SE})$, $\delta_k(z_{NW})$, $\delta_k(z_{NE})$ in case (iii).

Now, if $u^0 \in C_0(\Omega^*_0)$, we define $u^0_{k+1} \in C_0(\Omega^*_0)$ by first specifying the sets $\Omega^*_0(u^0_{k+1}, c)$, $1 \leq \mu \leq 2$, according to

$$
\Omega^*_1 \circ (u^0_{k+1}, c) = \{ z \in \Omega^*_0 \cap N^2_{k+1}(z) \cap \Omega^*_0 \},
$$

and then setting

$$
u^0_{k+1}(z) = \begin{cases} 
    (\beta^k v_k)(z), & \text{if } z \in \Omega^*_0 \circ (u^0_{k+1}, c) \\
    \psi_{k+1}(z), & \text{if } z \in \Omega^*_0 \circ (u^0_{k+1}, c) \\
    0, & \text{if } z \in \Gamma^*_k+1,0 \\
    d_i, & \text{if } z \in \Gamma^*_k+1,i, \quad 1 \leq i \leq m.
\end{cases}
$$

Prespecifying accuracy bounds $\varepsilon^l > 0$, $\mu = 1, 2$, on all levels $0 \leq k \leq l$, the nested iteration process combined with the scheme MGEPT results in

**ALGORITHM NMGEPT** ($u_{l,c}, (\varepsilon^l)_{k=0}^l, (\varepsilon^l)_{k=0}^l, C$):

Step 1. Set $k = 0$, choose $u^0_{0,c} \in C_0(\Omega^*_0)$ by $u^0_{0,c}(z) = \psi_0(z)$, $z \in \Omega^*_0$, and compute $u^0_{0,c}$ by performing $\text{MGEPT}(u^0_{0,c}, (\varepsilon^0)_{k=0}^l, (\varepsilon^l)_{k=0}^l, C)$.

Step 2. Set $k = k + 1$, determine $u^0_{k,c}$ by (48) and compute $u^0_{k,c}$ by performing $\text{MGEPT}(u^0_{k,c}, (\varepsilon^l)_{k=0}^l, (\varepsilon^l)_{k=0}^l, C)$.

Step 3. If $k = l$, stop the algorithm, otherwise go to Step 2.
We remark that in general the interpolated start iterate $u^0_{i,c}$ is not a supersolution to (23) on level $k$, but due to the global convergence of the scheme EPT this does not affect the convergence of the nested iteration scheme.

5. NUMERICAL RESULTS

As test examples we have chosen cylinders with a different number $m$ of cavities having circular or rectangular cross sections $\Omega$ and $\Omega_i$ respectively, $1 \leq i \leq m$, and being subjected to a torsion of varying twist angle per unit length $C$.

In particular, we have considered the following cases,

I. Circular cross sections

\[ \Omega = \{(x_1, x_2) \mid (x_1 - 0.5)^2 + (x_2 - 0.5)^2 < 0.25\} \]
\[ \Omega_i = \{(x_1, x_2) \mid (x_1 - x_i^1)^2 + (x_2 - x_i^2)^2 < r_i^2\}, \quad 1 \leq i \leq m \]

(i) $m = 0$ [cf. Figures 1, 5]

(ii) $m = 2$ [cf. Figures 3, 8]

\[ x_1^i = x_2^i = 0.5, \quad x_1^2 = 0.825, \quad r_1 = r_2 = 0.125 \]

(iii) $m = 9$ [cf. Figures 4, 11]

\[ x_1^i = x_2^i = 0.5, \quad r_1 = 0.2 \]
\[ x_1^2 = 0.15, \quad x_2^2 = 0.5; \quad x_1^3 = x_2^3 = 0.25 \]
\[ x_1^4 = 0.25, \quad x_2^4 = 0.75; \quad x_1^5 = 0.5, \quad x_2^5 = 0.15 \]
\[ x_1^6 = 0.5, \quad x_2^6 = 0.85; \quad x_1^7 = 0.75, \quad x_2^7 = 0.25 \]
\[ x_1^8 = 0.75; \quad x_1^9 = 0.85, \quad x_2^9 = 0.5 \]
\[ r_i = 0.05, \quad 2 \leq i \leq 9 \]

II. Rectangular cross sections

\[ \Omega = \{(x_1, x_2) \mid 0 < x_1 < 1, \quad 0 < x_2 < 1\}, \]
\[ \Omega_i = \{(x_1, x_2) \mid a_i^1 < x_1 < b_i^1, \quad a_i^2 < x_2 < b_i^2\}, \quad 1 \leq i \leq m \]

(i) $m = 0$ [cf. Figures 2, 6]

(ii) $m = 1$ [cf. Figure 7]

\[ a_1^1 = 0.2, \quad b_1^1 = 0.5, \quad a_1^2 = 0.65, \quad b_1^2 = 0.95 \]

(iii) $m = 2$ [cf. Figure 9]

\[ a_1^1 = 0.2, \quad b_1^1 = 0.4, \quad a_1^2 = 0.2, \quad b_1^2 = 0.8 \]
\[ a_2^1 = 0.6, \quad b_2^1 = 0.8, \quad a_2^2 = 0.2, \quad b_2^2 = 0.8 \]

(iv) $m = 4$ [cf. Figure 10]

\[ a_1^1 = 0.2, \quad b_1^1 = 0.4, \quad a_1^2 = 0.2, \quad b_1^2 = 0.4 \]
\[ a_2^1 = 0.6, \quad b_2^1 = 0.8, \quad a_2^2 = 0.2, \quad b_2^2 = 0.4 \]
For discretization of the corresponding LC-problems (12) we have chosen equidistant grids of step-sizes \( h_{k+1} = h_k/2 \), \( 0 \leq k \leq 4 \), with \( h_0 = 0.5 \).

In all cases the numerical results† clearly indicated the superiority of NMGEPT compared to MGEPT, thus illustrating the benefits of nested iteration. Indeed, computing a start-iterate on the finest grid by nested iteration, in all test examples it took at most three iterations until the final configuration of the elastic/plastic region was attained, while for this, using MGEPT with \( u^0 = \psi \) as start-iterate, considerably more iterations had to be performed (except for higher values of \( C \)).

We have also tested various configurations of the multi-grid cycles, in particular the cases \( \gamma_k = 1, 0 \leq k \leq l \) ("V-cycle") and \( \gamma_k = 2, 0 \leq k \leq l \) ("W-cycle"), with smoothing only before as well as before and after the defect correction process.

As a measure for the performance of the schemes we have computed the residues

\[
 r_i = \left[ \left( \sum_{x \in \Omega^i} |(\delta^i u_i)(x) + 2C|^2 \right) / N_i \right]^{1/2}
\]

with respect to the inactive grid-points.

Table I shows the computed residues obtained by NMGEPT in case of test example I. (ii) \( (C = 5.0, l = 5) \) for V-cycles and W-cycles with two smoothing iterations before or before and after the defect correction, \( \nu \) counting the iterations on the finest grid. For computation of \( u_5^0 \) we have chosen \( \nu = 1 \) iteration on level \( k = 0 \) and \( \nu = 2 \) iterations on levels \( 1 \leq k \leq 4 \) using \( \delta_k = \delta_k^2 = 1.0 \times 10^{-3} \), \( 0 \leq k \leq 4 \), as termination criterion. In all cases the final distribution of active/inactive grid-points was reached after three iterations on level \( l = 5 \) (this explains the relatively poor convergence for \( \nu = 3 \) and, in particular, the deterioration of \( r_5 \) from \( \nu = 1 \) to \( \nu = 2 \) in column 3 of Table I).

It is clear that the application of multi-grid techniques remains meaningful as long as \( k_{min} < l \), while it does not make sense if the degree of plastification is so high in relation to the grid-size \( h_i \) that no inactive grid-point can be found on the next coarser grid \( \Omega_{l-1} \). Indeed, in this case MGEPT simply reduces to scheme EPT with respect to the fixed grid \( \Omega_l \).

For test example I. (ii) Table II gives the values of \( k_{min} \) for \( l = 4, 5, 6 \) and torsion angles (per unit length) \( C = 2.5, 5.0 \) and 10.0.

<table>
<thead>
<tr>
<th>Table I. Example I. (ii), ( C = 5.0, l = 5 )</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \gamma_k = 1 ), ( 0 \leq k \leq 5 )</td>
</tr>
<tr>
<td>( \nu )</td>
</tr>
<tr>
<td>1</td>
</tr>
<tr>
<td>2</td>
</tr>
<tr>
<td>3</td>
</tr>
<tr>
<td>4</td>
</tr>
<tr>
<td>5</td>
</tr>
</tbody>
</table>

†All reported results in this section have been obtained by computations being performed on the Cyber 175 of ZRZ, TU Berlin.
Table II. Example I. (ii)

<table>
<thead>
<tr>
<th>$k_{\min}$</th>
<th>$C = 2.5$</th>
<th>$C = 5.0$</th>
<th>$C = 10.0$</th>
</tr>
</thead>
<tbody>
<tr>
<td>4</td>
<td>3</td>
<td>3</td>
<td>4</td>
</tr>
<tr>
<td>5</td>
<td>3</td>
<td>4</td>
<td>4</td>
</tr>
<tr>
<td>6</td>
<td>3</td>
<td>4</td>
<td>5</td>
</tr>
</tbody>
</table>

In Figures 1–4 we have represented the dependence of the plastification for various test examples on the applied torsion where active grid-points characterizing the discrete plastic region are marked by points.

Figures 5–11 show the plastification, the equipotential lines which correspond to the stress trajectories and the graph of the stress potential.

Note that the graph of the stress potential has a meaning in terms of Prandtl’s membrane analogon and Nadai’s sand-hill analogon which have been used in mechanical engineering for
experimental modelling of the torsion problem. First, using transparent material a ‘roof’ is constructed representing the graph of the generalized distance function. Second, a membrane is put across the boundary of the cross section and finally, a pressure of magnitude \(2C\) is applied to the membrane from opposite to the roof. Then, the deformation of the membrane corresponds to the stress potential, and the projection onto the plane of the boundary of the contact zones between membrane and roof indicates the free boundary between the elastic and the plastic region.

All plots shown in Figures 1–10 are based on numerical results computed on level \(l = 5\) by means of the multi-grid algorithm NMGEPT.

REFERENCES