Complex Boundary Elements for Contaminant Transport

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ABSTRACT

Recent numerical advances in the Complex Variable Boundary Element Method (CVBEM) provide easier-to-use analysis procedures in the study of advection-dominated contaminant transport of conservative species migration in a steady groundwater flow field. In this paper, the CVBEM is applied to groundwater advection contaminant transport problems, and the CVBEM numerical error is evaluated by use of the approximate boundary graphical technique. Besides providing actual solutions to many groundwater flow and advective contaminant transport problems, the CVBEM can be used to develop analytic test cases to be used in numerically calibrating other groundwater and contaminant transport numerical models for other classes of problems. In this paper, the CVBEM model is developed by use of a coupled $L^2$/Collocation fit to prescribed boundary conditions.

Key-_words: Complex Variables; Numerical Modeling; Contaminant Transport; Groundwater; CVBEM; Complex Variable Boundary Element Method

PURPOSE OF PAPER

The purpose of this paper is to report on the advances made in using the Complex Variable Boundary Element Method (or CVBEM) in the approximation of conservative species contaminant transport problems. Specifically, the CVBEM is a numerical boundary integral equation technique that is useful in describing flow characteristics of solute transport (for conservative species transport that are transported by groundwater flow hydrayics (advection)) such as flow paths and steady groundwater flow timing properties. In this paper, the use of the CVBEM in contaminant transport problems is extended by employment of a coupled least-squares optimization fit ($L^2$) and collocation fit in meeting the problem's prescribed boundary conditions. Because the CVBEM utilizes approximation functions that exactly solve the Laplace equation, there is no error in satisfying the governing potential flow linear operator equation except in meeting prescribed boundary conditions. The coupled $L^2$/Collocation error approach results in another technique to minimizing this approximation error. The "approximate boundary" associated to the CVBEM solution (i.e., the locus of points where the CVBEM solution achieves the prescribed boundary condition values) is used to graphically demonstrate the accuracy of the CVBEM model in achieving the solution to the groundwater flow problem.

MATHEMATICAL MODEL DEVELOPMENT

Several publications develop the CVBEM in depth, and the reader is directed to Hromadka and Lai (1987) for a thorough

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treatment of the numerical technique including FORTRAN computer codes to implement the technique. Publications that focus upon use of the CVBEM in contaminant transport problems include Hromadka and Yen (1986) and Yen and Hromadka (1986). In all of these prior works, the CVBEM is based upon a collocation technique in meeting boundary condition values. Use of an $L^2$ norm with the CVBEM can be found in Hromadka and others (1987). The current paper is the first report of a coupled $L^2$/Collocation method in matching boundary conditions by the CVBEM. A major advantage of the $L^2$/Collocation method over just $L^2$ or Collocation alone, is that the CVBEM "approximate boundary" (see Hromadka and LaI, 1987) is more readily developed (primarily at the interface between prescribed groundwater flow potential and normal flux boundary conditions), and the departures (i.e., distance) between the approximate and true problem boundaries are diminished over use of the Collocation method alone.

**COMPLEX VARIABLE BOUNDARY ELEMENTS (CVBEM)**

The CVBEM is a numerical approximation of the Cauchy integral formula of an analytic function $\omega(z)$ over a domain $\Omega$ and simple closed contour $\Gamma$ within $\Omega$,

$$\omega(z) = \frac{1}{2\pi i} \int_{\Gamma} \frac{\omega(\zeta) d\zeta}{\zeta - z} \quad (1)$$

where $\omega(z) = \phi(z) + i\psi(z)$; $\phi(z)$ is the two-dimensional potential function and $\psi(z)$ is the two-dimensional conjugate of $\phi(z)$; and $\omega(z)$ are known on $\Gamma$.

Flux values of the potential, $\phi(z)$, on $\Gamma$ are related to $\psi(z)$ by the Cauchy-Riemann relationship

$$\begin{vmatrix} \partial\phi \\ \partial n \end{vmatrix} = -i \begin{vmatrix} \partial\psi \\ \partial s \end{vmatrix} \quad (2)$$

where $n$ and $s$ are normal and tangential vector components along $\Gamma$, respectively.

Equation (1) is approximately solved by use of (for example) one-dimensional spline functions of $\phi(z)$ and $\psi(z)$ defined on small segments (boundary elements, $\Gamma_j$) of $\Gamma$. Nodal points, with coordinates $z_j$, $j=1,2,\ldots,n$, are placed on $\Gamma$, and each boundary element $\Gamma_j$ is a straight-line segment with endpoints $z_j$ and $z_{j+1}$. A global approximation function of $\omega(z)$ on the union of boundary elements, $\bigcup \Gamma_j$, is defined by $G_1(z)$ where

$$G_1(z) = \sum_{j=1}^{n} N_j(z) \omega_j \quad (3)$$

where $\omega_j = \phi_j + i\psi_j$ are nodal point $j$ values of $\omega(z_j)$, $\phi(z_j)$, $\psi(z_j)$; and $N_j(z)$ is a piecewise linear complex polynomial satisfying for all $z \in \bigcup \Gamma_j$:

$$N_j(z) = \begin{cases} 1, & z = z_j \\ 0, & z = z_{j-1}, z_{j+1} \\ 0, & z \notin \Gamma_{j-1} \cup \Gamma_{j} \end{cases} \quad (4)$$

Substituting the $G_1(z)$ approximation (the subscript 1 refers to a first-order spline function family) into the Cauchy integral equation results in

$$\hat{\omega}_1(z) = \frac{1}{2\pi i} \int_{\Gamma} \frac{G_1(z) \, d\zeta}{\zeta - z} \quad (5)$$

where $\hat{\omega}_1(z)$ is the CVBEM approximation, which is a linear combination of $n$ pairs of nodal point values $\phi_j$ and $\psi_j$, wherein $n$ values are known via prescribed boundary conditions, leaving no more than $n$ unknown nodal point values. Convergence of $\hat{\omega}_1(z)$ to the exact solution of the potential flow problem (i.e., steady state groundwater or advective contaminant transport) depends on the uniform convergence of $G_1(z)$ to $\omega(z)$ on $\Gamma$.

The aforementioned references present two techniques in developing estimates for the unknown CVBEM nodal values by use of a collocation technique, and by use of a $L^2$ technique. The coupled $L^2$/Collocation technique is described in the following section of this paper.

$L^2$/COLLOCATION

The CVBEM approximation function, $\hat{\omega}_1(z)$, based upon first order spline functions, can be written as (for problems
unsolved by a first order complex polynomial

\[ \hat{\varphi}_1(z) = \sum_{j=1}^{n} C_j(z - z_j) \ln(z - z_j) \tag{6} \]

where \( \ln \) is the complex natural logarithm function and \( C_j \) are complex constants \( C_j = a_j + i b_j \) for \( j = 1, 2, \ldots, n \). A collocation technique to solving for the CVBEM unknown constants (i.e., unknown nodal values) is to set implicit relationships (Hromadka and Lai, 1987)

\[
\begin{align*}
\text{Im}(\hat{\varphi}(z_j)) &= \psi(z_j) \\
\text{Re}(\hat{\varphi}(z_j)) &= \phi(z_j)
\end{align*}
\tag{7}
\]

where \( \text{Im} \) and \( \text{Re} \) are the imaginary and real components of \( \hat{\varphi}(z_j) \), respectively, and \( \psi(z_j) \) or \( \phi(z_j) \) are the unknown nodal values, respectively. Such an implicit relationship results in a Fredholm equation of the second kind.

An \( L^2 \) relationship is to minimize the integral objective function, \( \mathcal{X} \),

\[
\mathcal{X} = \int_{\Gamma_\phi} (\hat{\phi} - \phi)^2 \, d\Gamma + \int_{\Gamma_\psi} \left( \frac{\partial \hat{\psi}}{\partial n} - \frac{\partial \psi}{\partial n} \right)^2 \, d\Gamma \tag{8}
\]

where \( \Gamma_\phi \) and \( \Gamma_\psi \) refer to the prescribed potential and normal flux boundary conditions, respectively. The numerical procedure for minimizing \( \mathcal{X} \) can be found by the usual Gram-Schmidt procedures as discussed in Hromadka and others (1987).

The coupled \( L^2 / \text{Collocation} \) procedure is a four step process:

1. Locate all "interface" nodal points on \( \Gamma \) wherein both \( \phi \) and the normal derivative of \( \psi \) are known.
2. At all interface nodal points, \( z_k \), set dual nodal equations:
   \[
   \begin{align*}
   \text{Re}(\hat{\varphi}(z_k)) &= \phi(z_k) = \phi_k \\
   \text{Im}(\hat{\varphi}(z_k)) &= \psi(z_k) = \psi_k
   \end{align*}
   \]
   where both \( \phi_k \) and \( \psi_k \) are known nodal values.
3. At all non-interface nodal points, solve for unknown nodal values by the \( L^2 \) formulation (excluding the interface nodal points already solved).
4. An additional fourth step is used for computational purposes in developing the CVBEM "approximate boundary" (see next section):

   At all boundary corners (or locations in \( \bigcup \Gamma \) where \( \Gamma_j \) and \( \Gamma_j \) join at an angle \( \phi \pi \) at \( z_j \)), collocate the known boundary condition (say \( \phi \) is known at \( z_j \) to equate to \( \hat{\varphi}(z_j) \):
   \[
   \text{Re}(\hat{\varphi}(z_j)) = \phi_j
   \]
   or similarly,
   \[
   \text{Im}(\hat{\varphi}(z_j)) = \psi_j
   \]
   where \( \phi_j \) (or \( \psi_j \)) is known at \( z_j \).

**APPENDIX BOUNDARY**

The CVBEM approximation function \( \hat{\varphi}(z) = \phi(z) + i \psi(z) \) has the property that both \( \hat{\varphi}(z) \) and \( \hat{\psi}(z) \) exactly solve the Laplace equation in the interior of \( \Omega \), both are continuous functions on \( \Gamma \), and both are conjugate functions. Additionally, depending on the success in the numerical model by placement and number of nodal points on \( \Gamma \) (and hence the goodness of fit of \( G_j(z,x) \) to \( \omega(z) \) on \( \Gamma \)), \( \hat{\phi}(z) = \phi(z) \) on \( \Gamma_\phi \) and \( \hat{\psi}(z) = \psi(z) \) on \( \Gamma_\psi \). The "approximate boundary" concept graphically demonstrates the success in \( \hat{\phi}(z) \) matching the prescribed boundary condition values by the development of the approximate boundary, \( \Gamma \), which is the locus of points where \( \hat{\phi}(z) \) achieves the boundary condition values. (Hromadka and Lai (1987) develop the approximate boundary concept in detail.)

From steps (2) and (4) of the previous section, one is guaranteed that \( \Gamma \) and \( \Gamma \) coincide at all "interface" nodal points and all corners of \( \bigcup \Gamma \) (and hence corners of \( \Gamma \)). Excluding corners of \( \bigcup \Gamma \), let \( z^* \) be a point on \( \Gamma \) (which is a straight-line segment in \( \epsilon \) neighborhood of \( z^* \)). Then the approximate
boundary is located with normal and
tangential coordinates \((n, s) = (n^*, s^*)\) such that 
\(z^* = (0, s^*)\) and \(|n^*|\) is the normal
distance between \(\bar{\Gamma}\) and \(\Gamma\) with 
\(s = s^*\). \(\bar{\Gamma}\) is
found by locating where \(\hat{\phi}(n^*, s^*) = \phi(0, s^*)\)
on \(\Gamma_\bar{\phi}\) and where \(\psi(n^*, s^*) = \psi(0, s^*)\) on 
\(\Gamma_\psi\) (or where \(\frac{\partial \psi}{\partial n}(n^*, s^*) = \frac{\partial \psi}{\partial n}(0, s^*)\) on 
\(\Gamma_\psi\)). It is
apparent that at all \(n^*\) interface
nodes, \(\hat{\omega}(0, s^*) = \omega(0, s^*)\), and hence \(\hat{\Gamma}\) and \(\Gamma\)
intersect at interface nodes, by step (3) of
the \(L^2\)/Collocation procedure.

Using a dense set of \((0, s^*)\) point
locations on \(\Gamma\), the approximate boundary \(\hat{\Gamma}\)
is found, that corresponds to the CVBEM
\(\hat{\omega}(z)\), and the goodness of the CVBEM
approximation is readily seen by comparing
\(\hat{\Gamma}\) and \(\Gamma\). Should \(\hat{\Gamma} = \Gamma\) everywhere, then
\(\hat{\omega}(z) = \omega(z)\) on \(\Omega\). Additionally, if the analyst
is willing to accept \(\hat{\Gamma}\) as the "true"
boundary of the problem, then \(\hat{\omega}(z)\) is the "true"
solution of the redefined boundary value
problem.

**SOURCES AND SINKS**

The CVBEM develops an analytic
function on \(\Omega\). Sources and sinks are
introduced by adding the complex logarithm
function \(\ln(z - z_F)\) (Yen and Hromadka, 1986)
for each source sink at location \(z_F\), to the
CVBEM approximation function, \(\omega(z)\). The
composite CVBEM approximation is analytic
everywhere in \(\Omega\) except for all locations \(z_F\),
and the branch-cut originating from each \(z_F\).

**APPLICATION TO ADVECTIVE
CONTAMINANT TRANSPORT PROBLEMS**

In order to demonstrate the above
described procedures three selected
contaminant transport problems are
considered, the first two which demonstrate
the accuracy achieved as manifested by the
closeness between the approximate and true
boundaries of the problem. In all problems
considered, equally spaced nodes are used.
Other nodal placements are possible, with
similar results.

**Source-Sink Problem**

In this illustrative problem, a simple
groundwater well and injection well are
installed in a flow-field. Figures 1a,b,c
demonstrate the approximate boundary \(\hat{\Gamma}\)
developed from use of CVBEM approximations for two different nodal point
configurations. In Figure 1c, the closeness
of \(\hat{\Gamma}\) to \(\Gamma\) suggests that the underlying
CVBEM approximation function is a good
estimator of the background potential

**Landfill Evaluation**

Figure 2 depicts the application of the
CVBEM and approximate boundary
techniques to a landfill contaminant
transport problem. Injection wells are
problem is further studied regarding contaminant arrival times.

Application A. Here we will consider the steady flow produced by a single pumping well (50 m³/hr) near a landfill site with an equipotential boundary (Φ = 2 m) along the coordinate y = 1000 m. As shown in Figure 3, it takes the contaminant front produced by the landfill 9.0 years to reach the pumping well.

Figure 2.
(a) Problem Definition,
(b) CVBEM approximate boundary for 60 nodes,
(c) CVBEM approximate boundary for 70 nodes.

c onsidered to offset the effects of contaminant flow to the potable water. In Figures 2b and 2c, the addition of CVBEM nodal points improves the modeling accuracy as manifested by the reduction in departure between the approximate boundary, \( \Gamma \), and the true boundary, \( \Gamma^* \). The maximum departure of 40 m between \( \Gamma \) and \( \Gamma^* \) in Figure 2c is negligible in comparison to the overall scale of the problem; that is, the magnitude of difference between energy head values along \( \Gamma \) are minor.

Contaminant Transport Arrival Time

In the following two applications, the CVBEM landfill contaminant transport
Application B. When two injection wells are installed between the landfill and the pumping well, their influence on retarding the contaminant movement can be assessed. When 10 m³/hr is injected at each well it takes more than 13 years for the contaminant front to reach the pumping well (see Figure 4).

The results of Figures 3 and 4 are based upon a 75-node CVBEM model. Additionally, the approximate boundaries \( \Gamma \) for each CVBEM model deviated only a few meters from the true problem boundary shape, \( \Gamma \), and from the geometry of the landfill boundary. That is, \( \tilde{\Gamma} \) deviated from \( \Gamma \) well within the measuring tolerance possible in the field. Hence, the CVBEM results are exact should \( \tilde{\Gamma} \) be considered the true geometry of the problem. The \( L^2/\text{Collocation} \) approach was useful in these applications due to the reduction in effort needed to develop the approximate boundaries for numerical error evaluation and subsequent CVBEM model fine-tuning.

CONCLUSIONS

The CVBEM is applied to groundwater advection contaminant transport problems, and the CVBEM numerical error is evaluated by use of the approximate boundary graphical technique. Besides providing actual solutions to many groundwater flow and advective contaminant transport problems, the CVBEM can be used to develop analytic test cases to be used in numerically calibrating other groundwater and contaminant transport numerical models for other classes of problems. In this paper, the CVBEM model is developed by use of a coupled \( L^2/\text{Collocation} \) fit to prescribed boundary conditions.

The coupled \( L^2/\text{Collocation} \) method used in the CVBEM results presented herein was found to be useful in the development of the approximate boundary. The \( L^2/\text{Collocation} \) method was similar in computational effort to either the \( L^2 \) or Collocation approaches separately, but significantly simplified the plotting and definition of approximate boundaries corresponding to the CVBEM models.

REFERENCES


