

# Real and Complex Variable Models of Potential Flow

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Comparison of Real Variable and Complex Variable Approaches to Models of Potential Flow

# Section 1

## Methodologies

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# Real and Complex Variable Models

## Method of Fundamental Solutions

$$g_{rv}(\mathbf{x}, \mathbf{x}') = -\frac{1}{2\pi} \log |\mathbf{x} - \mathbf{x}'|$$

$$\hat{\phi}_{rv}(\mathbf{x}) = \sum_{j=1}^n a_j \log |\mathbf{x} - \mathbf{x}'_j|, \quad a_j \in \mathbb{R}$$

$$\text{where } \mathbf{x} = \begin{bmatrix} x \\ y \end{bmatrix}$$

## Complex Variable Simple Pole Approach

$$g_{cv}(z, z') = \frac{1}{z - z'}$$

$$\hat{\phi}_{cv}(\mathbf{x}) = \Re \left[ \sum_{j=1}^n \frac{c_j}{z - z_j} \right], \quad c_j \in \mathbb{C}$$

where  $z = x + iy$

# The General CVBEM Approximation Function

- The CVBEM approximation function is a linear combination of complex variable functions that are **analytic** within a given problem domain,  $\Omega$ :

$$\hat{\omega}(z) = \sum_{j=1}^n c_j g_j(z), \quad z \in \Omega, \quad (1)$$

- where

- ▶  $c_j = \alpha_j + i\beta_j$  are complex coefficients,
- ▶  $g_j(z)$  are analytic complex variable basis functions,
- ▶  $n$  is the number of basis functions being used in the approximation

- In the collocation approach, each term in the approximation function corresponds to **one** node and **two** collocation points.

Some possible  
basis functions:

- $(z - z_j) \ln_{\alpha_j}(z - z_j)$
- $1/(z - z_j)$
- $(z - z_j)^j$
- Digamma
- Polygamma
- And more!

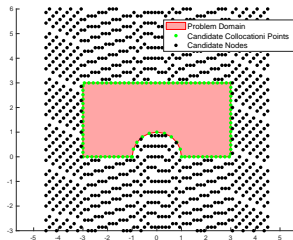
Hromadka II, T.V., Guymon, G.L., A Complex Variable Boundary Element Method: Development. *International Journal for Numerical Methods in Engineering*, pp. 25-37, 1984.

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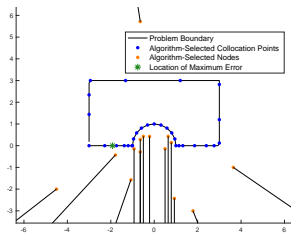
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# Node Position Algorithm (NPA)



**Figure:** Candidate nodes are **tested one-at-a-time** to determine which node contributes the most to reducing the maximum error of the MFS or CVBEM approximation function.

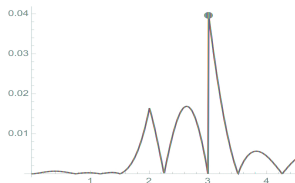
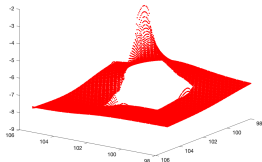


**Figure:** After a node is selected, the maximum error of the resulting CVBEM model is assessed. Two new collocation points are added at the two highest local maxima of the error function.

Demoes, N.J., Bann, G.T., Wilkins, B.D., Grubaugh, K.E. & Hromadka II, T.V., Optimization Algorithm for Locating Computational Nodal Points in the Method of Fundamental Solutions to Improve Computational Accuracy in Geosciences Modeling. *The Professional Geologist*, pp. 6-12, 2019.

# Error Computation

- The **maximum principle of harmonic functions** states that the maximum value of a harmonic function restricted to a particular domain occurs on the boundary of that domain.
- Let  $\Omega$  denote the problem domain. Our basis functions are constructed to be harmonic in  $\Omega$ . The target function is harmonic in  $\Omega$ . Therefore,  $\phi - \hat{\phi}$  is harmonic in  $\Omega$ .
- Thus,  $\max_{(x,y) \in \bar{\Omega}} |\phi - \hat{\phi}| = \max_{(x,y) \in \partial\Omega} |\phi - \hat{\phi}|$



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# Handling Mixed Boundary Conditions

The Dirichlet boundary conditions:

$$\begin{aligned}\hat{\phi}(x_{i,D}, y_{i,D}) &= \sum_{j=1}^n \alpha_j \lambda_j(x_{i,D}, y_{i,D}) - \beta_j \mu_j(x_{i,D}, y_{i,D}) \\ &= \phi(x_{i,D}, y_{i,D}), \\ \text{for } i &= 1, \dots, N_D, \quad (x_{i,D}, y_{i,D}) \in \partial\Omega_D.\end{aligned}$$

Let:

$$\begin{aligned}x_D &= \begin{bmatrix} x_{1,D} \\ x_{2,D} \\ \vdots \\ x_{N_D,D} \end{bmatrix}, \quad y_D = \begin{bmatrix} y_{1,D} \\ y_{2,D} \\ \vdots \\ y_{N_D,D} \end{bmatrix}, \\ x_N &= \begin{bmatrix} x_{1,N} \\ x_{2,N} \\ \vdots \\ x_{N_N,N} \end{bmatrix}, \quad \text{and } y_N = \begin{bmatrix} y_{1,N} \\ y_{2,N} \\ \vdots \\ y_{N_N,N} \end{bmatrix}.\end{aligned}$$

The Neumann boundary conditions:

$$\begin{aligned}\hat{\psi}(x_{i,N}, y_{i,N}) &= \sum_{j=1}^n \alpha_j \mu_j(x_{i,N}, y_{i,N}) + \beta_j \lambda_j(x_{i,N}, y_{i,N}) \\ &= \text{const}, \\ \text{for } i &= 1, \dots, N_N, \quad (x_{i,N}, y_{i,N}) \in \partial\Omega_N.\end{aligned}$$

In matrix form:

$$\begin{bmatrix} \underbrace{\phi(x_D, y_D)}_{N_D \times 1} \\ \underbrace{f}_{N_N \times 1} \end{bmatrix} = \begin{bmatrix} \underbrace{\lambda(x_D, y_D)}_{N_D \times n} & \underbrace{-\mu(x_D, y_D)}_{N_D \times n} \\ \underbrace{\mu(x_N, y_N)}_{N_N \times n} & \underbrace{\lambda(x_N, y_N)}_{N_N \times n} \end{bmatrix} \begin{bmatrix} \underbrace{\alpha}_{n \times 1} \\ \underbrace{\beta}_{n \times 1} \end{bmatrix},$$

where  $f = \gamma \begin{bmatrix} 1 \\ 1 \\ 1 \\ \vdots \\ 1 \end{bmatrix}$ ,  $\gamma \in \mathbb{R}$ .

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## Section 2

### Example Problems and Results

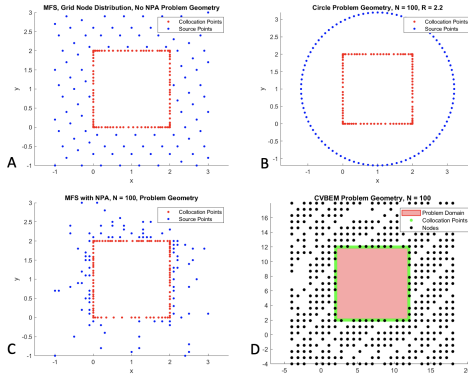
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# Corner Problem: Four Approaches

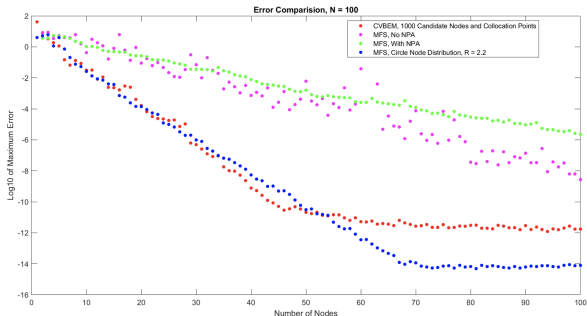


**Figure:** Problem geometry plots for the four different approaches. For all approaches, 100 nodes and collocations were used.

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# Error Comparison



**Figure:** Comparison of the maximum error plots from CVBEM method and three different MFS methods. CVBEM outperforms the grid node distribution and MFS with the NPA, but the circle node distribution with optimized radius is able to outperform the CVBEM.

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# Additional Results

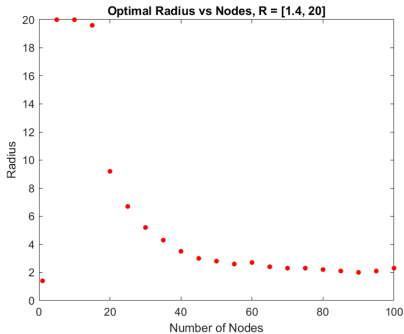


Figure: Optimal circle radius for up to  $n = 100$  nodes.

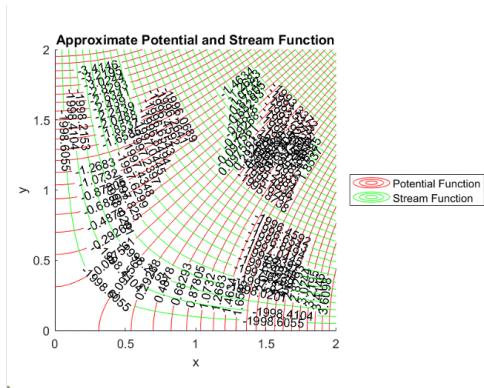


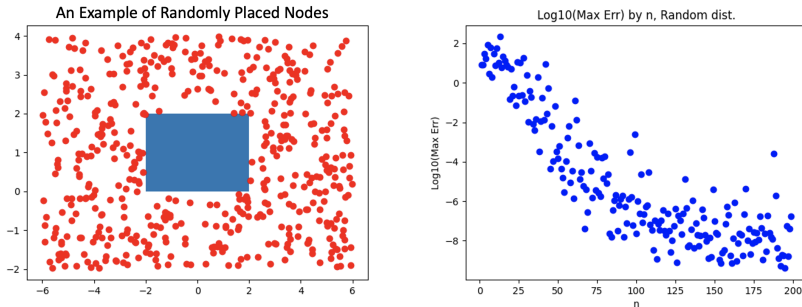
Figure: Approximate potential and stream functions for  $n = 50$  nodes.

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# NPA with Gaussian Randomness

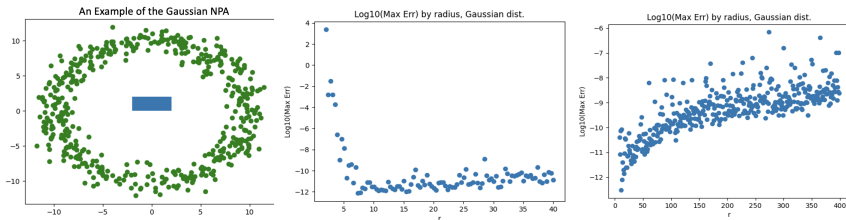


**Figure:** (Left) Sample distribution of candidate nodes with random placement. (Right)  $\log_{10}$  of maximum error for various-sized MFS models.

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# NPA with Gaussian Randomness - Radius Examination



**Figure:** (Left) Sample distribution of candidate nodes at a specified radius with Gaussian randomness applied. (Center)  $\log_{10}$  of maximum error for an MFS model as the radius of nodes is varied. Radii between 1 and 40 were examined. (Right)  $\log_{10}$  of maximum error for an MFS model as the radius of nodes is varied. Radii between 1 and 400 were examined.

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## Section 3

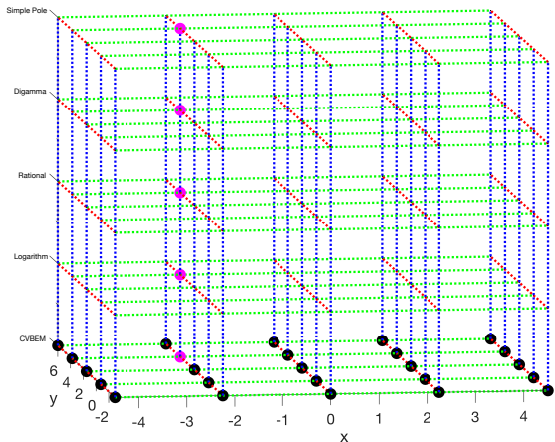
### Final Thoughts - Upcoming Research

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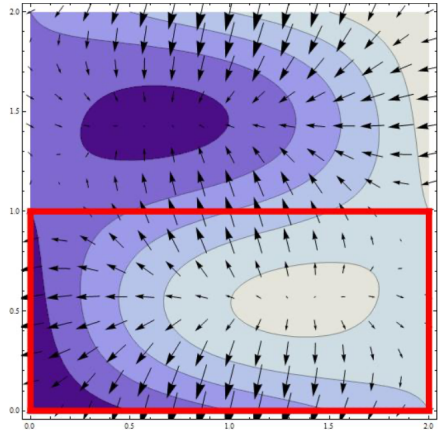
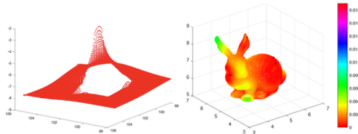
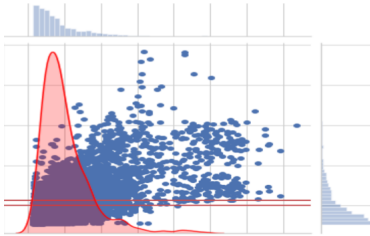
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# Composite Approximation Function with Line-search NPA



**Figure:** An alternative to the existing NPA is this algorithm, which proposes determining the location of each node, and then determining which basis function from a set of candidate functions to use at that particular location.

# Questions



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