

Analyzing layered soil problems with an alternative multi-region BEM technique and a new infinite boundary element formulation

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Abstract

The main objective of this work is to obtain an efficient three-dimensional boundary element (BE) formulation for layered soil simulation. This formulation is obtained by combining an alternative multi-region technique with an infinite boundary element (IBE) formulation. Kelvin fundamental solutions are employed, considering the static analysis of isotropic and linear-elastic domains.

Establishing relations between the displacement fundamental solutions of the different domains, the alternative technique used in this paper allows analyzing all domains as a single solid, not requiring equilibrium or compatibility equations. This approach also leads to a smaller system of equations when compared to the usual subregion technique, and the results obtained are more accurate.

The two-dimensionally mapped infinite boundary element (IBE) formulation here used is based on a triangular BE with linear shape functions. One advantage of this formulation over quadratic or higher order elements is that no additional degrees of freedom are added to the original BE mesh by the presence of the IBEs. Thus, the IBEs allow the mesh to be reduced without compromising the accuracy of the result.

The use of IBEs improves the advantages of the alternative multi-region technique, contributing for the low computational cost and allowing a considerable mesh reduction. Furthermore, the results show good agreement with the ones given in other work, confirming the accuracy of the presented formulation.

Keywords: infinite boundary elements, alternative multi-region technique, three-dimensional soil, static analysis.



1 Introduction

Considering specifically infinite multi-domain models, many options are available in the literature and each one of them implies on advantages and disadvantages. However, depending on the problem to be solved, one technique may become more attractive than the others.

Most cases, a numerical approach may be employed. The finite element method (FEM) is still popular (Karakus *et al.* [1]), however has some disadvantages when compared to other options such as the boundary element method (BEM). The FEM requires the discretization of the infinite domain, implying on a high number of elements and leading to a large and sometimes impracticable processing time. To reduce these inconveniences, some authors use infinite elements together with finite elements, such as Sadecka [2].

It becomes more viable to solve these problems with the BEM, once only the boundary of the domains requires discretization. This allows reducing the problem dimension, implying on less processing time. This advantage is explored in several works (Almeida and Paiva [3]) and more developments are making the BEM even more attractive to future applications.

The classical way to consider domains in contact with the BEM, which is described in details by Brebbia and Dominguez [4], is based on imposing equilibrium and compatibility conditions for all interface points between every pair of domains in contact. These impositions may cause inaccuracies, and numerous blocks of zeros are generated at the final system of equations.

However, Ribeiro and Paiva [5] present an alternative multi-region BEM technique, which does not require equilibrium nor compatibility conditions along the interfaces. Considering a constant Poisson ratio, it is possible to establish relations between the displacement fundamental solutions and to analyze all subdomains as a single solid. Thus, a better continuity between domains in contact is guaranteed and therefore the result accuracy is improved. In addition to that, no blocks of zeros are present at the final system of equations, which is reduced. Thus, better results are obtained in less processing time.

Another way to improve the BEM performance is by using infinite boundary elements (IBEs). The first reference to an IBE was Kagawa *et al.* [6], in which the shape functions of an origin BE are multiplied by special decay functions. Another type of IBE may be obtained by using mapped functions to relate the local system of coordinates to the global one, as originally shown by Beer and Watson [7]. In those studies that make use of two-dimensional IBEs, such as performed by Moser *et al.* [8], it may be noted that they are generally based on quadrilateral BEs. An alternative to these type of IBE is given by Ribeiro and Paiva [9], which presented a mapped IBE based on a triangular BE with linear shape functions. One advantage of this approach over quadratic or higher order elements is that no additional degrees of freedom are added to the original BE mesh by the presence of the IBEs.

The aim of this work is to combine the two BE techniques presented in [5, 9], obtaining a new and more efficient multi-region BE formulation for layered soil



simulation. The soil is modeled with variable elasticity module and a constant Poisson ratio, as described by Gibson [10]. The results obtained are consistent with those of another author, confirming the accuracy of the presented approach. In addition to that, the use of IBEs contributed for the low computational cost, allowing a considerable mesh reduction.

2 Boundary element formulation

The equilibrium of a solid body can be represented by a boundary integral equation called the Somigliana Identity, which for homogeneous, isotropic and linear-elastic domains is

$$c_{ij}(y) u_j(y) + \int_{\Gamma} p_{ij}^*(x, y) u_j(x) d\Gamma(x) = \int_{\Gamma} u_{ij}^*(x, y) p_j(x) d\Gamma(x) \quad (1)$$

Eqn (1) is written for a source point y at the boundary, where the displacement is $u_j(y)$. The constant c_{ij} depends on the Poisson ratio and the boundary geometry at y . The field point x goes through the whole boundary Γ , where displacements are $u_j(x)$ and tractions are $p_j(x)$. The integral kernels $u_{ij}^*(x, y)$ and $p_{ij}^*(x, y)$ are Kelvin three-dimensional fundamental solutions for displacements and tractions, respectively. Kernel $u_{ij}^*(x, y)$ has order $1/r$ and kernel $p_{ij}^*(x, y)$ order $1/r^2$, where $r = |x - y|$, so the integrals have singularity problems when x approaches y . Therefore the stronger singular integral, over the traction kernel, has to be defined in terms of a Cauchy Principal Value (CPV).

To solve eqn (1) numerically, the boundary is divided into subregions within which displacements and tractions are approximated by known shape functions. Here these subregions are of two types, finite boundary elements (BEs) and infinite boundary elements (IBEs). The BEs employed are triangular, as shown in fig. 1 with the local system of coordinates, $\xi_1 \xi_2$, and the local node numbering. The

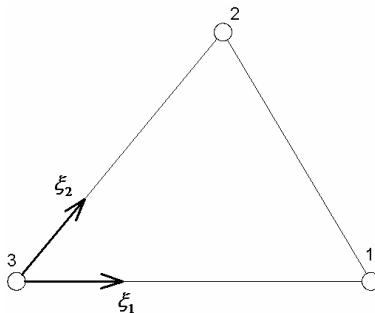


Figure 1: Triangular boundary element.

following approximations are used for this BE:

$$u_j = \sum_{k=1}^3 N^k u_j^k, \quad p_j = \sum_{k=1}^3 N^k p_j^k \quad (2)$$

Eqn (2) relate the boundary values u_j and p_j to the nodal values of the BE. The BEs have 3 nodes and for each node there are three components of displacement u_j^k and traction p_j^k . The shape functions N^k used for these approximations are

$$N^1 = \xi_1, \quad N^2 = \xi_2, \quad N^3 = 1 - \xi_1 - \xi_2 \quad (3)$$

The same shape functions are used to approximate the boundary geometry:

$$x_j = \sum_{k=1}^3 N^k x_j^k \quad (4)$$

where x_j^k are the node coordinates. The same functions are also used to interpolate displacements and tractions for the IBEs:

$$u_j = \sum_{k=1}^{Np} N^k u_j^k, \quad p_j = \sum_{k=1}^{Np} N^k p_j^k \quad (5)$$

Each IBE has Np nodes and not the 3 that the BEs have. The IBE geometry, on the other hand, is approximated by special mapping functions, as discussed in more detail in Section 3.

By substituting eqns (2) and (5) in eqn (1), eqn (6) is obtained:

$$c_{ij}(y) u_j(y) + \sum_{e=1}^{N_{BE}} \left\{ \sum_{k=1}^3 [\Delta p_{ij}^{ek} u_j^k] \right\} + \sum_{e=1}^{N_{IBE}} \left\{ \sum_{k=1}^{Np} [\Delta^\infty p_{ij}^{ek} u_j^k] \right\} = \sum_{e=1}^{N_{BE}} \left\{ \sum_{k=1}^3 [\Delta u_{ij}^{ek} p_j^k] \right\} + \sum_{e=1}^{N_{IBE}} \left\{ \sum_{k=1}^{Np} [\Delta^\infty u_{ij}^{ek} p_j^k] \right\} \quad (6)$$

N_{BE} is the number of BEs and N_{IBE} is the number of IBEs. For BEs:

$$\Delta p_{ij}^{ek} = \int_{\gamma_e} |J| N^k p_{ij}^* (x, y) d\gamma_e, \quad \Delta u_{ij}^{ek} = \int_{\gamma_e} |J| N^k u_{ij}^* (x, y) d\gamma_e \quad (7)$$

In eqn (7), γ_e represents the domain of element e in the local coordinate system and the global system of coordinates is transformed to the local one by the Jacobian $|J| = 2A$, where A is the element area in the global system. On the other hand, for IBEs:

$$\Delta^\infty p_{ij}^{ek} = \int_{\gamma_e} |\infty J| N^k p_{ij}^* (x, y) d\gamma_e, \quad \Delta^\infty u_{ij}^{ek} = \int_{\gamma_e} |\infty J| N^k u_{ij}^* (x, y) d\gamma_e \quad (8)$$

Eqn (8) is analogous to eqn (7), and the calculation of Jacobian $|\infty J|$ is discussed in Section 3. Integrals of eqns (7) and (8) are calculated by standard



BEM techniques. Non-singular integrals are evaluated numerically by using integration points. The singular ones, on the other hand, are evaluated by the technique presented in reference [11]. Finally, the free term c_{ij} may be obtained by rigid body motions.

Writing eqn (6) for all boundary nodes leads to the following system:

$$\Delta p \cdot u = \Delta u \cdot p \tag{9}$$

The Δp_{ij}^{ek} and $\Delta^\infty p_{ij}^{ek}$ element contributions, including the free term c_{ij} , are assembled into matrix Δp , while Δu_{ij}^{ek} and $\Delta^\infty u_{ij}^{ek}$ contributions are assembled into matrix Δu . Vectors u and p contain all boundary displacements and tractions, respectively. Reorganizing this system so as to separate the known boundary values from the unknown yields a system of equations whose solution is all the unknown boundary values.

3 Infinite boundary elements

Three types of mapping are considered, as illustrated in fig. 2.

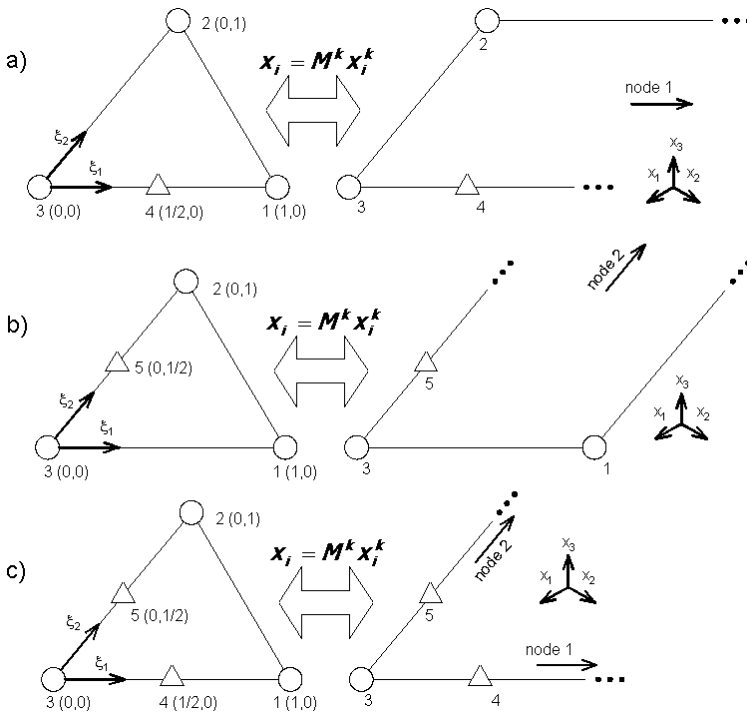


Figure 2: Types of mapping.

In the first type, represented in fig. 2a, only direction ξ_1 is mapped and node 1 is placed at infinity. The IBE is represented in the local coordinate system on the left and in the global coordinate system on the right. The global coordinates x_i are related to the local ones by special mapping functions, M^k , and the nodal global coordinates, x_i^k . Node 4 is created only to replace node 1 in the mapping and does not contribute to the integrals.

Fig. 2b is analogous to fig. 2a, but in this case only direction ξ_2 is mapped and node 2 is placed at infinity. Therefore, node 5 is created to facilitate the mapping. Finally, in fig. 2c both local directions are mapped and nodes 1 and 2 are placed at infinity. As a result, the auxiliary nodes 4 and 5 must be created to replace them in the mapping.

Ribeiro and Paiva [9] use auxiliary coordinates $\bar{\xi}_1$ and $\bar{\xi}_2$ to obtain the mapping functions for each case. When only direction ξ_1 is mapped, the result is:

$$M_{1\infty}^4 = \bar{\xi}_1(\xi_1) = \frac{\xi_1}{1 - \xi_1} \quad (10)$$

$$M_{1\infty}^2 = \xi_2 \quad (11)$$

$$M_{1\infty}^3 = 1 - \bar{\xi}_1(\xi_1) - \xi_2 = 1 - \frac{\xi_1}{1 - \xi_1} - \xi_2 \quad (12)$$

The symbol “ 1∞ ” is used to indicate that these expressions are valid if only direction ξ_1 is mapped. These functions are then employed to relate the local system of coordinates to the global one. In other words:

$$x_i = M_{1\infty}^4 x_i^4 + M_{1\infty}^2 x_i^2 + M_{1\infty}^3 x_i^3 \quad (13)$$

After obtaining eqn (13), the Jacobian used when only direction ξ_1 is mapped may be calculated as follows:

$$|{}^\infty J_1| = \frac{\partial x_1}{\partial \xi_1} \frac{\partial x_2}{\partial \xi_2} - \frac{\partial x_2}{\partial \xi_1} \frac{\partial x_1}{\partial \xi_2} = \frac{2A_1}{(1 - \xi_1)^2} \quad (14)$$

where A_1 is the area of the triangle drawn between nodes 2, 3 and 4 in the global system of coordinates.

For mapping only in direction ξ_2 , the functions obtained are:

$$M_{2\infty}^1 = \xi_1 \quad (15)$$

$$M_{2\infty}^5 = \bar{\xi}_2(\xi_2) = \frac{\xi_2}{1 - \xi_2} \quad (16)$$

$$M_{2\infty}^3 = 1 - \xi_1 - \bar{\xi}_2(\xi_2) = 1 - \xi_1 - \frac{\xi_2}{1 - \xi_2} \quad (17)$$

The symbol “ 2∞ ” is used to indicate that only direction ξ_2 is mapped. Therefore, the global system is related to the local one as follows:

$$x_i = M_{2\infty}^1 x_i^1 + M_{2\infty}^5 x_i^5 + M_{2\infty}^3 x_i^3 \quad (18)$$



and the Jacobian is

$$|{}^{\infty}J_2| = \frac{2A_2}{(1 - \xi_2)^2} \quad (19)$$

where A_2 refers to the area of the triangle drawn between nodes 1, 3 and 5 in the global system of coordinates.

Finally, for mapping in both directions ξ_1 and ξ_2 , the mapping functions are

$$M_{\infty}^4 = \frac{\xi_1}{1 - \xi_1} \quad (20)$$

$$M_{\infty}^5 = \frac{\xi_2}{1 - \xi_2} \quad (21)$$

$$M_{\infty}^3 = 1 - \frac{\xi_1}{1 - \xi_1} - \frac{\xi_2}{1 - \xi_2} \quad (22)$$

The symbol “ ∞ ” is used to indicate that both directions are mapped. The local system of coordinates is related to the global one as follows:

$$x_i = M_{\infty}^4 x_i^4 + M_{\infty}^5 x_i^5 + M_{\infty}^3 x_i^3 \quad (23)$$

and the Jacobian is now

$$|{}^{\infty}J_3| = \frac{2A_3}{(1 - \xi_1)^2 (1 - \xi_2)^2} \quad (24)$$

where A_3 is the area of the triangle drawn between nodes 3, 4 and 5 in the global system.

4 Alternative multi-region formulation

In fig. 3 a problem with two regions and submitted to arbitrary boundary conditions is presented.

The regions have the same Poisson ratio ν and different elasticity modules, E_1 for region Ω_1 and E_2 for region Ω_2 . The boundary Γ_1 of region Ω_1 is divided in two parts, Γ_{12} and $\bar{\Gamma}_1$. The part of Γ_1 which is in contact with region Ω_2 is denoted by Γ_{12} and the rest of it is named $\bar{\Gamma}_1$. Analogously, region Ω_2 boundary Γ_2 is divided into Γ_{21} for the contact and $\bar{\Gamma}_2$ for the free surface. Consequently:

$$\Gamma_1 = \bar{\Gamma}_1 + \Gamma_{12}, \quad \Gamma_2 = \bar{\Gamma}_2 + \Gamma_{21} \quad (25)$$

Kelvin displacement fundamental solutions for regions Ω_1 and Ω_2 may be written as:

$$u_{ij1}^* = \frac{1}{16\pi\mu_1(1 - \nu)r} [(3\nu - 4)\delta_{ij} + r_{,i}r_{,j}] \quad (26)$$

$$u_{ij2}^* = \frac{1}{16\pi\mu_2(1 - \nu)r} [(3\nu - 4)\delta_{ij} + r_{,i}r_{,j}] \quad (27)$$



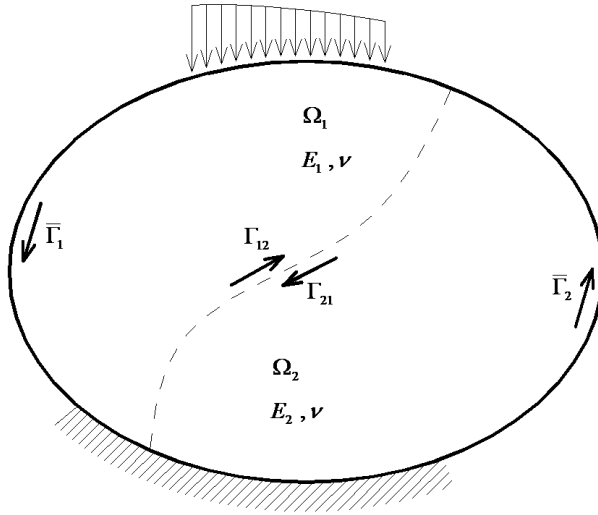


Figure 3: Problem with two regions.

where

$$\mu_1 = \frac{E_1}{2(1 + \nu)} \tag{28}$$

$$\mu_2 = \frac{E_2}{2(1 + \nu)} \tag{29}$$

In such a way, the fundamental solutions may be related as follows:

$$u_{ij2}^* = \frac{E_1}{E_2} u_{ij1}^* \tag{30}$$

Eqn (30) may be rewritten as:

$$u_{ij2}^* = u_{ij1}^* + \frac{\Delta E_{12}}{E_1} u_{ij2}^* \tag{31}$$

where:

$$\Delta E_{ij} = E_i - E_j \tag{32}$$

Kelvin traction fundamental solution may be written as

$$p_{ij}^* = \frac{-1}{8\pi(1 - \nu)r^2} \left[\frac{\partial r}{\partial \eta} [(1 - 2\nu)\delta_{ij} + 3r_{,i}r_{,j}] + (1 - 2\nu)(\eta_i r_{,j} - \eta_j r_{,i}) \right] \tag{33}$$

One may observe that eqn (33) do not depend on the elasticity module, than it is the same for domains Ω_1 and Ω_2 . Thus, Kelvin traction fundamental solution may be represented as p_{ij}^* for both domains.

The objective is to obtain a single integral equation which considers both regions. In order to do that, the first step is to write the Somigliana Identity integrating only the boundary of region Ω_1 . Maintaining the same source point, the Somigliana Identity is then again written but now integrating only the boundary of region Ω_2 . By adding these two expressions, an initial version of the final equation is obtained. The next step is to substitute eqn (31) into this initial version and to develop the expression. This demonstration is detailed by Ribeiro and Paiva [5], and the final result is

$$\begin{aligned} \left[c_{ij1} + \frac{E_2}{E_1} c_{ij2} \right] u_j + \int_{\bar{\Gamma}_1} p_{ij}^* u_j d\bar{\Gamma}_1 + \frac{E_2}{E_1} \int_{\bar{\Gamma}_2} p_{ij}^* u_j d\bar{\Gamma}_2 - \frac{\Delta E_{12}}{E_1} \int_{\Gamma_{21}} p_{ij}^* u_j d\Gamma_{21} \\ = \int_{\bar{\Gamma}_1} u_{ij1}^* p_j d\bar{\Gamma}_1 + \int_{\bar{\Gamma}_2} u_{ij1}^* p_j d\bar{\Gamma}_2 \end{aligned} \quad (34)$$

Variables x and y were omitted in order to reduce the expression. Extending it to an arbitrary number of domains, it becomes:

$$\begin{aligned} \left\{ \sum_{s=1}^{nd} \left[\frac{E_s}{E_1} c_{ijs} \right] \right\} u_j + \sum_{e=1}^{ne} \left[\frac{E_e}{E_1} \int_{\bar{\Gamma}_e} p_{ij}^* u_j d\bar{\Gamma}_e \right] + \sum_{c=1}^{nc} \left[\frac{\Delta E_{mn}}{E_1} \int_{\Gamma_{mn}} p_{ij}^* u_j d\Gamma_{mn} \right] = \\ = \sum_{e=1}^{ne} \left[\int_{\bar{\Gamma}_e} u_{ij1}^* p_j d\bar{\Gamma}_e \right] \end{aligned} \quad (35)$$

In eqn (35), the total number of domains is nd , the number of contact boundaries is nc and the number of external boundaries is ne . The first summation represents the coefficient c_{ij} (y) of eqn (1), which contributes in the matrix Δp of eqn (9). Before calculating it all coefficients c_{ijs} , one for each domain, must be known. More details about this formulation are provided by Ribeiro and Paiva [5].

In Section 2, eqn (1) was used as a starting point to obtain the BEM system of equations which solution is the unknown boundary values. If the same steps are repeated for eqn (35), valid for multi-regions, a similar system of equations is obtained. The unknowns of this new system are the non-prescribed boundary values plus the interface displacements. The total number of unknowns is reduced when compared to the classic multi-region technique described in reference [4], once the interface tractions are not included in this case. This justifies why the alternative formulation leads to less time processing. A better interface continuity is also guaranteed, once all regions are modeled as a single solid.

5 Example

This example aims to analyze an infinite non-homogeneous domain problem with the proposed formulation. The domain considered is composed by two layers of different elasticity module, as illustrated in fig. 4.

Layer 1 has a 9000 kN/m^2 elasticity module, a 0.5 Poisson ratio and 15 m of thickness. Layer 2 has a 900 kN/m^2 elasticity module, a 0.5 Poisson ratio and



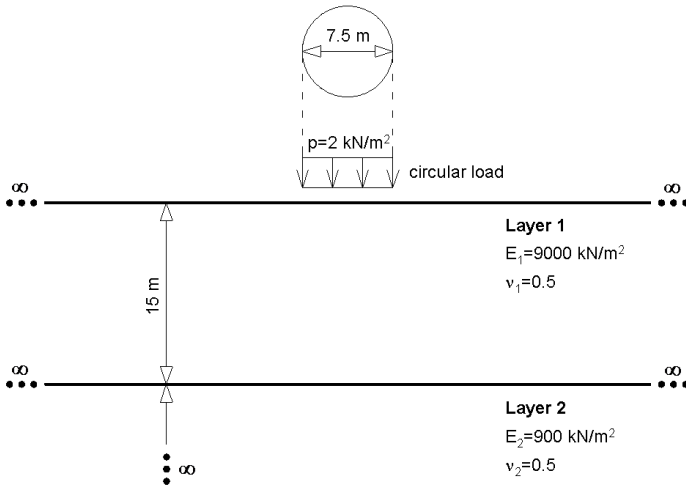


Figure 4: Layered soil problem.

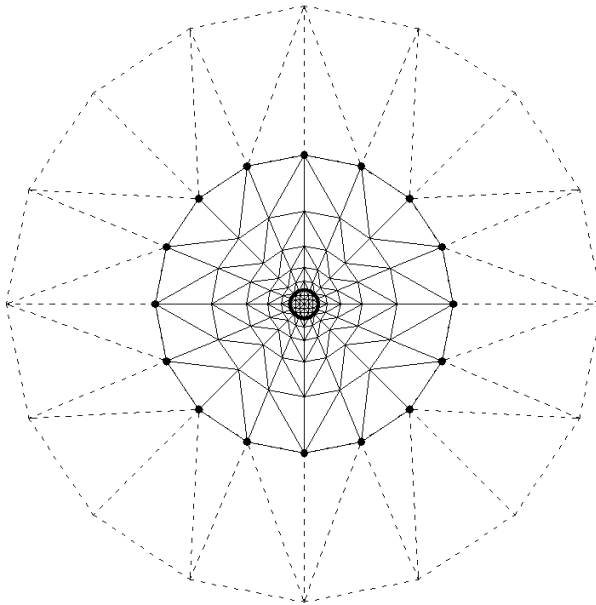


Figure 5: Mesh employed.

infinite thickness. Both layers are infinite in radial directions. A vertical circular 2 kN/m^2 uniform loading with a 7.5 m diameter is applied at the top layer surface.

This problem was simulated using two identical meshes, one for the surface and another for the contact between layers, totalizing 242 nodes, 448 BEs and 64 IBEs. The surface mesh may be visualized in fig. 5, in which the polygon outlined in the

Table 1: Vertical displacement ($m \times 10^{-3}$).

Number of nodes	No IBEs used	Error (%)	Mesh with IBEs	Error (%)
Analytical	2.5000	—————	2.5000	—————
242	2.0874	16.5	2.5123	0.5
274	2.3865	4.5	2.5209	0.8
306	2.5033	0.1	2.5212	0.8
338	2.5207	0.8	2.5213	0.9

center corresponds to the loaded area, the dashed lines represent the IBEs and the rest of the mesh is composed of BEs. The points marked at the perimeter of the BE mesh receive all the influence of the IBEs.

Using the values adopted in this example and applying the analytical solution given by Burmister [12], a $2.5000 \times 10^{-3} m$ vertical displacement is obtained for the central point of the circle. Simulating this problem with the mesh of fig. 5, a vertical displacement of $2.5123 \times 10^{-3} m$ was obtained. This value agrees with the analytical solution, with an error of 0.5%. In order to evaluate the influence of the IBEs, the example was simulated with the same BE mesh but without the IBEs. A displacement of $2.0874 \times 10^{-3} m$ was then obtained, with the higher error of 16.5% relative to the analytical value.

In order to improve this precision more BEs and degrees of freedom were added at the mesh limits. In this way, the values in table 1 were obtained.

As may be observed, 306 nodes were needed for the BE mesh to overcome the precision of the 242 node mesh with IBEs. Comparing these two values it may be concluded that, in this example, to maintain the error below 0.5%, the use of IBEs allows a mesh reduction of 21%.

6 Conclusions

In this paper, an alternative multi-region BE technique was combined with a IBE formulation in order to obtain an efficient numerical tool for layered soil simulation. Electing one domain as a reference and establishing relations between its displacement fundamental solution and the ones of the other regions allows integrating all domains as a single solid. This approach eliminates the need of equilibrium and compatibility relations between the different media. Therefore, better results may be obtained in less processing time.

The IBEs employed have the advantage of not increasing the original number of degrees of freedom, as demonstrated in the example presented. The results obtained with the IBEs showed good agreement with an analytical solution, and the use of IBEs promoted a mesh reduction of 21%.



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