

Volume Integral Equation Based Diakoptic Method for Electromagnetic Modeling

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Abstract—A novel diakoptic method based on volume integral equation (VIE) modeling of subsystems is proposed for 3-D electromagnetic analysis. The theoretical foundation of the method are the surface and volume equivalence principles, as it combines the VIE and surface integral equation (SIE) formulations, in conjunction with the method of moments (MoM). The method breaks the original structure into a number of non-overlapping closed-region subsystems that contain inhomogeneous dielectric materials and an open-region subsystem. Each subsystem is analyzed completely independently applying the double-higher-order large-domain Galerkin generalized MoM-VIE-SIE (VSIE) or MoM-SIE solvers. The final solution is obtained from diakoptic matrices expressing linear relations between the electric and magnetic equivalent surface current coefficients on diakoptic surfaces. The proposed VSIE-diakoptic method is validated, evaluated, and discussed in several characteristic examples. The examples demonstrate that the diakoptic method substantially increases the efficiency of the conventional MoM-VIE approach. When compared to the pure MoM-VIE double-higher-order technique, the diakoptic approach enables very considerable accelerations and memory savings, while fully preserving the accuracy of the analysis.

Index Terms—Numerical techniques, electromagnetic analysis, diakoptic analysis, volume integral equation, surface equivalence principle, method of moments, domain decomposition, higher order modeling.

I. INTRODUCTION

THE method of moments (MoM) in conjunction with the volume integral equation (VIE) formulation is a well-established approach to computational electromagnetic (CEM) modeling of three-dimensional (3-D) dielectric structures, as well as composite dielectric and metallic structures [1]. The

MoM-VIE methodology is especially suitable in cases involving inhomogeneous and complex dielectric materials. In such cases, it is a natural choice and is preferred over the overall more frequently used MoM methodology based on the surface integral equation (SIE) approach [2]. In many other cases, the MoM-VIE approach provides a useful alternative to the MoM-SIE modeling and on par or sometimes a more efficient solution. In fact, there has recently been a great renewed interest in VIE modeling in CEM and its applications, e.g., [3]-[5]. However, any increase in computational efficiency, in terms of either the computation time or the memory requirements, while preserving accuracy and robustness of this theoretically and conceptually very simple and elegant volumetric modeling technology, is always desired, especially if it can be done in a systematic, methodological fashion.

This paper proposes a novel diakoptic method based on VIE modeling of subsystems, as one possible strategy aimed at substantially enhancing the efficiency of the conventional MoM-VIE approach. The theoretical foundation of the method consists of the surface and volume equivalence principles. The numerical foundation of the method is a generalized MoM-VIE-SIE (VSIE) solver [6]-[9]. The diakoptic approach [10]-[14] solves a large and complex electromagnetic (EM) system as a linear combination of solutions of diakoptic subsystems, by means of matrix linear relations between coefficients in equivalent electric and magnetic surface current expansions on subsystem boundary surfaces (diakoptic surfaces).

Being based on the surface equivalence principle, the proposed diakoptic analysis is similar in spirit to the equivalence principle algorithm (EPA) [15]-[18] and the nested equivalence principle algorithm (NEPAL) [19], [20]. However, unlike in EPA, where the solution is based on establishing relations between the entire set of equivalent electric/magnetic currents of each domain with all the currents in all other domains and incident fields, matrix relations in diakoptics are established between electric and magnetic currents within each individual domain. Computation of these relations is solely based on manipulations with the current expansion coefficients. As such, it can be applied to the generalized MoM-VSIE method regardless of the geometry modeling (straight or curvilinear) and it can be implemented in already existing generalized MoM-VSIE codes with virtually no changes in those codes, except the partitioning. A difference is also that EPA methods introduce another set of

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basis functions for connecting the regions, while the diakoptic approach works with the same basis functions as the MoM-SIE. While NEPAL solves a VIE, it is based solely on the representation of the scattered field by radiation of point sources, represented by multipoles, and it is not obvious if and how it could be applied to more general or arbitrary VIE elements and basis functions. On the other hand, the general diakoptic analysis can certainly be seen as belonging, in a broader sense, to the class of contemporary domain decomposition (DD) methods, including those based on the finite element method (FEM) [21]-[27] and those based on the finite difference frequency-domain and time-domain (FDFD and FDTD) methods [28], [29], as well as DD-SIE methods [30], [31].

In specific, the VSIE-diakoptic method divides the original structure into a number of non-overlapping closed-region subsystems, which generally contain inhomogeneous dielectric materials and are analyzed by the generalized MoM-VSIE solver, and an open-region subsystem that encloses the other subsystems and is analyzed using the MoM-SIE technique. Parts of the original problem are represented by matrix linear relations, and diakoptic coefficients are formally similar to voltages and currents at ports of a linear multiport network. Diakoptic surfaces may completely coincide with the SIE surfaces, and this paper explicitly considers only such situations in numerical examples. However, in other situations, SIE surfaces may be inside diakoptic surfaces, where they may be discretized by the pure MoM-SIE method, which would introduce non-diakoptic surface-current unknowns in the VSIE-diakoptic model. In all cases, the unknowns on diakoptic surfaces are the only unknowns in the final system of equations. Closed-region diakoptic subsystems in the model can, but do not need to, be the same. The implemented VSIE-diakoptic numerical discretization is based on the double-higher-order large-domain Galerkin-type MoM-VIE and MoM-SIE modeling described in detail in [6] and [7], respectively. The hybridization of the VIE and SIE methods which results in a generalized volume-surface integral equation (VIE-SIE or VSIE) method is done as outlined in [8], [9]. The basic theory and preliminary results of the VSIE-diakoptic analysis are presented in a summary form in [14].

One more interesting interpretation of the diakoptic method is in order here. Namely, as the relation between the equivalent sources at the diakoptic surface is linear and includes contributions of the excitations situated in the subsystem, it can be written in a formally identical matrix form as the Norton (or Thévenin) representation of a linear multiport network. It is, therefore, valid for any EM environment (system) in which the subsystem may be placed (and combined with other subsystems of the same or different form), and the proposed diakoptic analysis can be formally interpreted as being equivalent to Norton's theorem in circuit theory. However, the diakoptic approach features the full EM-field generalization of the Norton representation, where, in fact, each pair of electric and magnetic equivalent surface current coefficients on the diakoptic surface represents a generalized port, to which a multiport network analogy is then

applied. So, the diakoptic approach is not a lumped-element approximation of EM fields, but rather a 3-D EM generalization, based on a rigorous Maxwell-type solver, of Norton's theorem.

The diakoptic method manipulates with linear relations between expansion coefficients for currents within a subsystem and equivalent currents at the diakoptic surface. Hence, results for electromagnetic subsystems once used in the diakoptic analysis can be stored for future reuse, with a minimum computational overhead. We can change one (or more) subsystems (e.g., change material composition, excitations, dimensions or placement of subcomponents, etc.) and rerun the analysis, to re-compute the diakoptic coefficients extremely expeditiously. For instance, if any part of the system is changed, only coefficients for that part have to be recalculated in the diakoptic model. This can be very beneficial to design and optimization procedures. Additionally, if the problem includes many identical subdomains, such as in analysis and design of antenna or scatterer arrays, the solution of one constitutive subsystem can be efficiently reused for congruent subsystems.

This paper has three principal goals. (1) It demonstrates that the proposed VSIE-diakoptic method substantially increases the efficiency of the conventional MoM-VIE approach, and this is the primary goal. When compared to the pure MoM-VIE technique implementing the same type of discretization (double-higher-order large-domain discretization in our case, i.e., [6]), the diakoptic approach enables very considerable accelerations and memory savings, while preserving the accuracy of the analysis. However, the paper does not claim that this is the only or the best strategy to enhance the efficiency of the conventional MoM-VIE method. (2) The paper also demonstrates that the general diakoptic methodology can effectively include VIE-based subsystems, and shows how that can be done. (3) Finally, the paper combines the general MoM-VIE methodology and the general DD methodology, which are two out of several general numerical methodologies in CEM. However, the paper does not speculate on comparisons in terms of capabilities and performance of the presented diakoptics-VIE approach with a possible contemporary DD-VIE approach.

The rest of the paper is organized as follows. Section II presents the theory and implementation of the novel VSIE-diakoptic method starting with the surface and volume equivalence principles and a set of VSIE equations for the unknown volume and surface equivalent currents. This also includes double-higher-order MoM-VSIE modeling of EM subsystems and diakoptic surfaces, derivation of linear relations between diakoptic coefficients, representation of EM subsystems by diakoptic matrices, and solution of the diakoptic matrix system. In Section III, the proposed VSIE-diakoptic method is validated, evaluated, and discussed in several characteristic examples.

II. DIAKOPTIC METHOD BASED ON MOM-VSIE MODELING

A. Surface and Volume Equivalence Principles

In the diakoptic analysis, the EM system is subdivided into

subsystems by virtue of the surface equivalence principle [32]. For the clarity of explanation, we consider a single domain v , bounded by a closed surface S , as depicted in Fig. 1(a), so two subsystems, v and its exterior. According to this principle, the electric and magnetic fields \mathbf{E} and \mathbf{H} everywhere can be uniquely evaluated by knowing electric and magnetic tangential fields \mathbf{E}_S and \mathbf{H}_S at S .

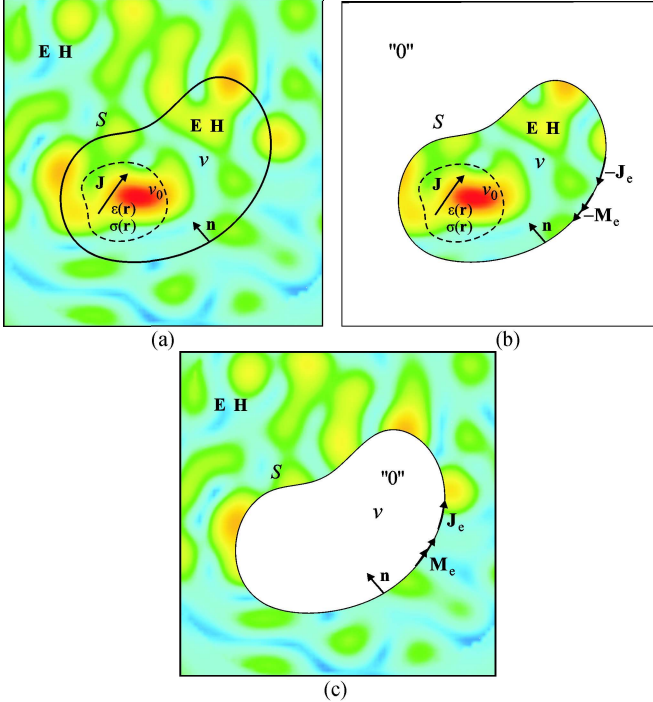


Fig. 1. Illustration of the surface and volume equivalence principles, as the theoretical foundation of the VSIE-diaoptical approach to electromagnetic analysis: (a) original EM system, divided into two subsystems by a closed surface S (diaoptical surface), on which equivalent surface currents are placed in order to radiate the actual fields, (b) equivalent interior problem, which includes an inhomogeneous dielectric body taken into account by the radiation of equivalent volume currents, and (c) equivalent exterior problem (with original fields outside and zero field inside S).

Let us assume that interior impressed volume electric currents of density \mathbf{J} exist only in a domain v_0 , which is a part of v . For example, these currents can represent an actual volume current distribution of a dielectric body located inside v . The fields \mathbf{E} and \mathbf{H} can be obtained as being radiated by fictitious electric and magnetic equivalent surface currents whose density vectors are given by $\mathbf{J}_e = -\mathbf{n} \times \mathbf{H}_S$ and $\mathbf{M}_e = \mathbf{n} \times \mathbf{E}_S$, where \mathbf{n} stands for a normal unit vector on S , directed toward the interior region, as shown in Fig. 1(a). As a result of subtraction of the equivalent currents and their fields from the original problem, the fields \mathbf{E} and \mathbf{H} in v remain unchanged, while the fields in the exterior region are annihilated, Fig. 1(b). The surface equivalence principle for the exterior region is established in the analogous way, Fig. 1(c).

Assume now that the domain v_0 in Fig. 1(a) actually represents an arbitrarily shaped inhomogeneous dielectric body, with the permittivity $\epsilon(\mathbf{r})$ and conductivity $\sigma(\mathbf{r})$ of the

dielectric material being known functions of position (i.e., the vector \mathbf{r}), while the permeability at all points is μ_0 . In accordance to the volume equivalence principle [32], the dielectric inhomogeneity can be taken into account by the radiation of volumetric electric currents of density \mathbf{J} , which we can consider as the impressed sources when applying the surface equivalence principle, in Fig. 1. The constitutive (material) relationship for the total electric field at each point of the domain v_0 and the tangential total electric and magnetic field boundary conditions on the surface S yield

$$\mathbf{E}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e) + \mathbf{E}^{\text{exc}} = \mathbf{D} / \epsilon_c \quad (1)$$

$$\mathbf{n} \times (\mathbf{E}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e) + \mathbf{E}^{\text{exc}}) = 0 \quad (2)$$

$$\mathbf{n} \times (\mathbf{H}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e) + \mathbf{H}^{\text{exc}}) = 0 \quad (3)$$

where $\mathbf{E}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e)$ and $\mathbf{H}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e)$ are field vectors due to volumetric electric current density and surface electric and magnetic currents, \mathbf{D} is the equivalent electric displacement (or flux density) vector, and ϵ_c is the equivalent complex permittivity of the material at that point, while \mathbf{E}^{exc} and \mathbf{H}^{exc} are excitation electric and magnetic fields resulting from arbitrarily positioned sources inside and/or outside of v . The vectors \mathbf{J} and \mathbf{D} are related to each other as

$$\mathbf{J} = j\omega C \mathbf{D}, \quad C = \frac{\epsilon_c - \epsilon_0}{\epsilon_c}, \quad \epsilon_c = \epsilon - j \frac{\sigma}{\omega}, \quad (4)$$

with ω being the angular frequency of \mathbf{E}^{exc} (time-harmonic convention $e^{j\omega t}$ is used) and C the electric contrast (with respect to the surrounding medium) at each point of the object. Given the integral expressions for the fields in terms of sources $\mathbf{E}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e)$ and $\mathbf{H}(\mathbf{D}, \mathbf{J}_e, \mathbf{M}_e)$ [6], [7], [14], (1)–(3) represent a system of three coupled electric/magnetic field integral VSIE equations with the sources \mathbf{D} , \mathbf{J}_e , and \mathbf{M}_e as unknown quantities; we discretize and solve this system using the MoM.

B. Double-Higher-Order MoM-VSIE Discretization

The VSIE-diaoptical method is based on the double-higher-order (in both geometrical and current approximation) MoM modeling. This facilitates utilization of large curved volume and surface elements. Volume elements are Lagrange-type curved hexahedra of geometrical orders K_u , K_v , and K_w ($K_u, K_v, K_w \geq 1$), shown in Fig. 2(a). Within each element, the vector \mathbf{D} is expanded via divergence-conforming hierarchical polynomial vector basis functions in parametric coordinates u , v , and w with arbitrary current-approximation orders N_u , N_v , and N_w ($N_u, N_v, N_w \geq 1$) [6], [14]. In the MoM-SIE discretization, modeling of S is carried out using Lagrange generalized curved quadrilateral elements of geometrical orders K_u and K_v ($K_u, K_v \geq 1$), in

Fig. 2(b), with the currents \mathbf{J}_e and \mathbf{M}_e over the patches being approximated by divergence-conforming hierarchical polynomial vector bases with orders N_u and N_v ($N_u, N_v \geq 1$) [7]. The K and N orders in one VIE or SIE element can be adopted completely independently from each other. All orders in different elements are also adopted at will, and there is no required connection between VIE and SIE orders. In the hybrid MoM-VSIE technique, the equations are tested using the Galerkin method.

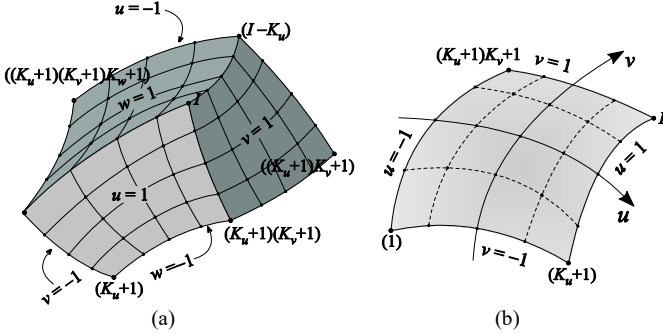


Fig. 2. Lagrange interpolation curved parametric elements for higher order VSIE-diaoptical analysis of the EM system in Fig. 1: (a) MoM-VIE generalized hexahedral element, determined by $I = (K_u + 1)(K_v + 1)(K_w + 1)$ interpolation nodes arbitrarily positioned in space, and (b) MoM-SIE generalized quadrilateral patch, defined by $I = (K_u + 1)(K_v + 1)$ nodes.

C. VSIE-Diaoptical Analysis

Discretizing (1)–(3) with the MoM, we obtain the following linear matrix system of equations:

$$\begin{array}{c} N_S \quad N_V \quad N_S \quad 1 \quad 1 \\ \leftrightarrow \quad \leftrightarrow \quad \leftrightarrow \quad \leftrightarrow \quad \leftrightarrow \\ N_S \quad \downarrow \left[\begin{array}{ccc} \mathbf{Z}_1 & \mathbf{Z}_2 & \mathbf{Z}_3 \\ \mathbf{Z}_4 & \mathbf{Z}_5 & \mathbf{Z}_6 \\ \mathbf{Z}_7 & \mathbf{Z}_8 & \mathbf{Z}_9 \end{array} \right] \left[\begin{array}{c} \mathbf{j}_e \\ \mathbf{d} \\ \mathbf{m}_e \end{array} \right] = \left[\begin{array}{c} \mathbf{v}_1 \\ \mathbf{v}_2 \\ \mathbf{v}_3 \end{array} \right] \\ N_V \quad \uparrow \\ N_S \quad \uparrow \end{array} \quad (5)$$

where $[\mathbf{d}]$, $[\mathbf{j}_e]$, and $[\mathbf{m}_e]$ are column-matrices of unknown coefficients that approximate \mathbf{D} , \mathbf{J}_e , and \mathbf{M}_e , respectively, the generalized MoM-VSIE system matrix, $[\mathbf{Z}]$, is partitioned into sub-matrices $[\mathbf{Z}_k]$, $k = 1, 2, \dots, 9$, and $[\mathbf{v}_l]$, $l = 1, 2, 3$, form the excitation column-matrix. The N_V equations in (5) in the middle, N_S equations at the top, and N_S equations at the bottom are obtained by discretizing (1), (2), and (3), respectively. In particular, sub-matrices $[\mathbf{Z}_5]$ and $[\mathbf{Z}_m]$, $m = 1, 3, 7, 9$ correspond to pure VIE and SIE discretizations, respectively, while $[\mathbf{Z}_n]$, $n = 2, 4, 6, 8$, result from the generalized VSIE analysis [8], [9], [14].

A combined field integral equation (CFIE) formulation for the diaoptical surface (surface S in Fig. 1) can be obtained by multiplying the first N_S equations in (5) by α , and the last N_S equations by β , where we adopt $\alpha = 1$ and $\beta = -\sqrt{\mu_0/\epsilon_0}$. The CFIE formulation is used to avoid numerical instabilities at

resonant frequencies of the diaoptical surface. Summing the new equations and moving all the terms associated with $[\mathbf{m}_e]$ to the right-hand side of the matrix equation results in

$$\begin{bmatrix} \alpha\mathbf{Z}_1 + \beta\mathbf{Z}_7 & \alpha\mathbf{Z}_2 + \beta\mathbf{Z}_8 \\ \mathbf{Z}_4 & \mathbf{Z}_5 \end{bmatrix} \begin{bmatrix} \mathbf{j}_e \\ \mathbf{d} \end{bmatrix} = - \begin{bmatrix} \alpha\mathbf{Z}_3 + \beta\mathbf{Z}_9 \\ \mathbf{Z}_6 \end{bmatrix} \begin{bmatrix} \mathbf{m}_e \\ \mathbf{v}_2 \end{bmatrix} + \begin{bmatrix} \alpha\mathbf{v}_1 + \beta\mathbf{v}_3 \\ \mathbf{v}_2 \end{bmatrix}, \quad (6)$$

which allows establishing the following matrix linear relation:

$$\begin{bmatrix} \mathbf{j}_e \\ \mathbf{d} \end{bmatrix} = \begin{bmatrix} \mathbf{Y} \\ \mathbf{C} \end{bmatrix} \begin{bmatrix} \mathbf{m}_e \\ \mathbf{d}_0 \end{bmatrix}. \quad (7)$$

Based on (7), each of the subsystems in Fig. 1, i.e., the interior subsystem, for which $i = 1$, and the exterior subsystem, with $i = 2$, can be described by linear relations between coefficients $[\mathbf{j}_e]$ and $[\mathbf{m}_e]$ cast in matrix form as

$$[\mathbf{j}_{ei}] = [\mathbf{Y}_i] [\mathbf{m}_{ei}] + [\mathbf{j}_{0i}], \quad i = 1, 2 \quad (8)$$

where $[\mathbf{Y}_i]$ is the $N_S \times N_S$ diaoptical matrix $[\mathbf{Y}]$ in (7) for the i -th subsystem and $[\mathbf{j}_{0i}]$ is the $N_S \times 1$ column-matrix that contains coefficients representing the excitation in the subsystem. These matrices are numerically calculated by considering the coefficients $[\mathbf{m}_{ei}]$ as excitation and coefficients $[\mathbf{j}_{ei}]$ as the response – in the i -th subsystem. To obtain $[\mathbf{Y}_i]$, we excite the subsystem with only one coefficient of $[\mathbf{m}_{ei}]$ set to unity value and all other coefficients (as well as all excitations within the subsystem) set to zero; with an excitation by the j -th coefficient in $[\mathbf{m}_{ei}]$, the coefficients of \mathbf{J}_e obtained using the generalized MoM-VSIE solver constitute the j -th column of $[\mathbf{Y}_i]$. From (7), we also obtain linear relations between $[\mathbf{m}_{e1}]$ and the VIE coefficients of \mathbf{D} ($i = 1$), given by

$$[\mathbf{d}] = [\mathbf{C}] [\mathbf{m}_{e1}] + [\mathbf{d}_0], \quad (9)$$

where the $N_V \times N_S$ diaoptical matrix $[\mathbf{C}]$ is filled column by column by computing the coefficients of \mathbf{D} in the same MoM-VSIE analysis with $[\mathbf{Y}_i]$. For example, the solution for $[\mathbf{j}_e]$ and $[\mathbf{d}]$ of the matrix equation

$$\begin{bmatrix} \alpha\mathbf{Z}_1 + \beta\mathbf{Z}_7 & \alpha\mathbf{Z}_2 + \beta\mathbf{Z}_8 \\ \mathbf{Z}_4 & \mathbf{Z}_5 \end{bmatrix} \begin{bmatrix} \mathbf{j}_e \\ \mathbf{d} \end{bmatrix} = - \begin{bmatrix} \alpha\mathbf{Z}_3 + \beta\mathbf{Z}_9 \\ \mathbf{Z}_6 \end{bmatrix} \begin{bmatrix} 1 \\ 0 \\ \vdots \\ 0 \end{bmatrix} \quad (10)$$

forms the first column of matrix $[\mathbf{Y} \mathbf{C}]^T$ in (7), and so on. In

order to obtain the column matrix $[\mathbf{j}_0 \mathbf{d}_0]^T$, we solve the above matrix equation with $[\alpha \mathbf{v}_1 + \beta \mathbf{v}_3 \mathbf{v}_2]^T$ on the right-hand side. While evaluating the diakoptic matrices, the matrix equation (6), i.e., the corresponding system of linear equations, is filled and LU (lower-upper) decomposed only once, and all the results are obtained with LU substitutions using different excitations. Namely, the main matrix on the left-hand side in (10) is LU factorized only once for multiple right-hand side vectors. Only forward/backward substitutions are applied in order to compute all the elements of the matrices $[\mathbf{Y}]$ and $[\mathbf{C}]$. The complexity of LU is $O(N^3)$ and the complexity of a pair of forward/backward substitution is $O(N^2)$. With this, the advantage of utilizing a direct solver in the diakoptic approach is straightforward. In addition, the proposed VSIE-diakoptic approach can be combined with fast direct solvers [33]-[35]. An iterative solver can also be applied to subsystems, and the calculation in subsystems can be accelerated by fast methods [36]-[38].

Note that, since the matrix relation in (8) relates the electric surface current to the magnetic surface current, it could be considered as a form of the generalized impedance boundary condition (GIBC) [39], [40].

The solution of the original EM problem, in Fig. 1(a), is obtained by relating the diakoptic coefficients in expansions of the electric and magnetic equivalent surface currents on S (on the diakoptic surfaces, i.e., individual subsystem boundaries). These relations come out to be as simple as

$$[\mathbf{j}_{e1}] = [\mathbf{j}_{e2}] = [\mathbf{j}_e] \text{ and } [\mathbf{m}_{e1}] = [\mathbf{m}_{e2}] = [\mathbf{m}_e] \quad (11)$$

where the use is made of the facts that the signs of equivalent sources are opposite (based on the equivalence theorem) and so are the normal unit vector directions for the two neighboring subsystems. Diakoptic boundary conditions in (11) lead to stable VSIE-diakoptic CFIE solutions in all cases considered and all tests performed. Note that, theoretically, (11) is directly derived from the boundary conditions for both fields \mathbf{H} and \mathbf{E} at the diakoptic surface. Combining (8) and (11), the following diakoptic matrix system is obtained:

$$([\mathbf{Y}_1] - [\mathbf{Y}_2])[\mathbf{m}_e] = -[\mathbf{j}_{01}] + [\mathbf{j}_{02}] \quad (12)$$

which is solved for $[\mathbf{m}_e]$ using a direct solver. Note that the matrices in (10) and (12) that need to be LU factorized (once) in the VSIE-diakoptic method are, in general, considerably smaller than the main matrix of the pure MoM-VIE technique implementing the same type of discretization. Once $[\mathbf{m}_e]$ is found, $[\mathbf{j}_e]$ and $[\mathbf{d}]$ can be computed from (8) and (9). Other quantities (e.g., electric and magnetic fields) can be calculated by post-processing of the coefficients by the MoM-VSIE code.

Note that when all the subsystems are isolated, there is one large exterior subsystem left to be solved by the MoM-SIE method. While this open-region subsystem relating all other subsystems becomes the bottleneck of the diakoptic method in

such cases, the complexity of the method, associated with this subsystem only, is still significantly reduced in comparison with the pure MoM-SIE approach. From a purely mathematical standpoint (omitting EM theoretical background of the method, which plays a key role in the diakoptic technique), using the diakoptic approach we solve two systems with N_S and one with $(N_S/M + N_V)$ numbers of unknowns, respectively, with M being the number of congruent subdomains, as shown in (6)-(12) and Appendix I, instead of solving a system with $2N_S$ unknowns. Therefore, in diakoptics, the linear size of the system of equations associated with the global exterior subsystem is halved. In addition, theoretically, any matrix solver that allows us to obtain (6), (7), (8), and (9), and to solve (12) can be used with the diakoptic approach.

Note also that transition from (5) to (6)-(7) can be considered as the first step of static condensation [41]. However, further steps of static condensation would require one to substitute vector $[\mathbf{j}_e \mathbf{d}]^T$ back into (3) and proceed further with its solution. In the presented diakoptic technique, however, matrix equation (5) is created for each of the subsystems constituting interior and exterior of the diakoptic surface in Fig. 1, resulting into two matrix equations that are solved simultaneously for shared set of unknowns. Depending on the subsystem, equations (5) may or may not contain unknown coefficients $[\mathbf{d}]$. Using the surface equivalence principle and the procedure of decomposing the problem into subsystems that are solved independently, while taking into account their interaction via equivalent surface currents, we obtain two relations between $[\mathbf{j}_e]$ and $[\mathbf{m}_e]$ as shown in (8), and then solve the final system in (12) as explained above.

The presented diakoptic analysis is generalized to the case of an arbitrary number of subsystems. Equations (11) are valid for a common diakoptic boundary of two neighboring subsystems and the matrix representations (8) and (9) stand for each subsystem, so the diakoptic system has the same form, since each part of a diakoptic surface is always shared by two adjacent subsystems. As an illustration, consider three subsystems, denoted as Ω_1 , Ω_2 , and Ω_3 , where the second and the third subsystem are in touch with the first, but not with each other. With the densities of the respective equivalent diakoptic current distributions being $(\mathbf{j}_{e1}, \mathbf{m}_{e1})$, $(\mathbf{j}_{e2}, \mathbf{m}_{e2})$, and $(\mathbf{j}_{e3}, \mathbf{m}_{e3})$, adjusting equations (11) and (12) leads to

$$\begin{bmatrix} \mathbf{j}_{e2} \\ \mathbf{j}_{e3} \end{bmatrix} = [\mathbf{j}_{e1}] = [\mathbf{j}_e] \text{ and } \begin{bmatrix} \mathbf{m}_{e2} \\ \mathbf{m}_{e3} \end{bmatrix} = [\mathbf{m}_{e1}] = [\mathbf{m}_e], \quad (13)$$

$$\left(\begin{bmatrix} \mathbf{Y}_2 \\ \mathbf{Y}_3 \end{bmatrix} - [\mathbf{Y}_1] \right) [\mathbf{m}_e] = - \begin{bmatrix} \mathbf{j}_{02} \\ \mathbf{j}_{03} \end{bmatrix} + [\mathbf{j}_{01}], \quad (14)$$

where matrices $[\mathbf{Y}_i]$ and $[\mathbf{j}_{0i}]$ correspond to the i -th subsystem ($i = 1, 2, 3$).

III. NUMERICAL EXAMPLES AND DISCUSSION

A. Array of Homogeneous Dielectric Cubical Scatterers

As the first example of the application and validation of the VSIE-diaoptics approach, we consider an 8×8 array of homogeneous lossless dielectric cubes of edge lengths $a = \lambda_0/6$, where λ_0 is the free-space wavelength, and relative permittivity $\epsilon_r = 2.25$, excited by a uniform plane electromagnetic wave from the direction described by $\theta_{\text{inc}} = 90^\circ$ and $\phi_{\text{inc}} = 0^\circ$, as shown in Fig. 3(a). The surface-to-surface distance between neighboring cubes is $d = \lambda_0/6$, in both x - and y -directions.

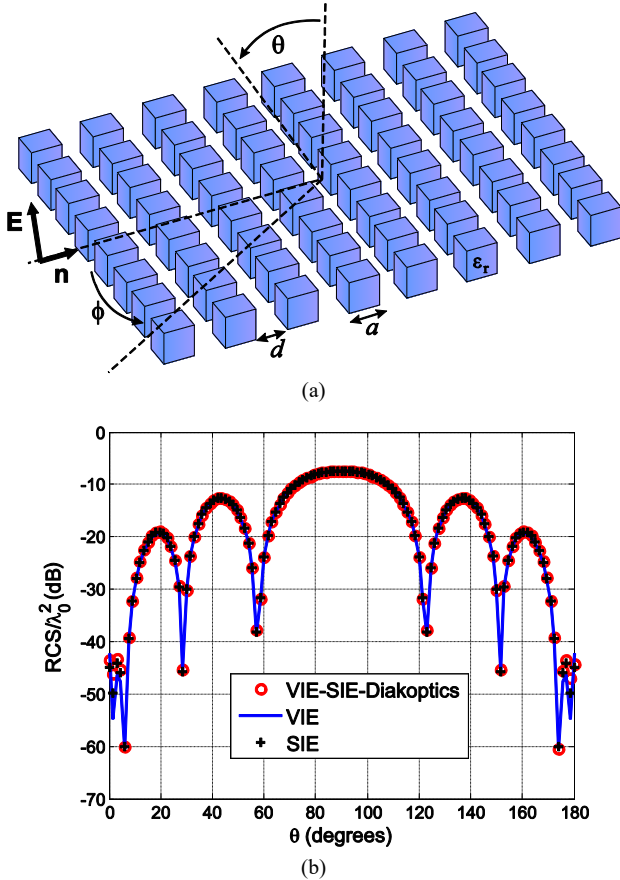


Fig. 3. (a) Array of homogeneous dielectric cubical scatterers excited by a uniform plane electromagnetic wave ($\theta_{\text{inc}} = 90^\circ$, $\phi_{\text{inc}} = 0^\circ$) and (b) normalized bistatic radar cross section (RCS) of the array in the $\phi = 0^\circ$ cut, computed by the proposed VSIE-diaoptics method, by the pure MoM-VIE technique [6], and the pure MoM-SIE technique [7].

For the diaoptics analysis, the system in Fig. 3(a) is split into 65 diaoptics subsystems: 64 interior VSIE subsystems coinciding with dielectric cubes and a free-space exterior SIE subsystem, which is in touch with all VSIE subsystems. A single VIE element is used for every cubical diaoptics subsystem, and the orders are adopted to be $K_u = K_v = K_w = 1$ and $N_u = N_v = N_w = 3$. Each cubical diaoptics surface is modeled by six SIE patches with $K_u = K_v = 1$ and $N_u = N_v = 2$. The number of VIE

unknowns per subsystem is $N_v = 108$ and the total number of diaoptics unknowns (summing the unknowns for all diaoptics surfaces) amounts to $2N_s = 6,144$, which results in a total of $N_{\text{tot}}^{\text{Dia}} = 6,252$ unknowns and $T_{\text{tot}}^{\text{Dia}} = 98$ s of computation time for the analysis. In all computations presented in this paper, no parallelization or EM system symmetries are used. All computations are performed on an Intel® Core™2 Quad CPU Q9550 at 2.83 GHz, with 8 GB RAM, under 64-bit Windows 7 operating system. Shown in Fig. 3(b) is the normalized bistatic radar cross section (RCS) of the array in a plane defined by $\phi = 0^\circ$. A comparison is made between the diaoptics solution and the results obtained by the pure MoM-VIE method (with LU decomposition) [6], which require $N_{\text{tot}}^{\text{VIE}} = 6,912$ unknowns and $T_{\text{tot}}^{\text{VIE}} = 439$ s of computation time, and an excellent agreement is observed. In this example, the advantage of the diaoptics approach is not so much in the reduction of the number of unknowns, but of the computation time (about 4.5 times). Moreover, the computer memory (RAM) consumption is $M_{\text{tot}}^{\text{Dia}} = 151$ MB for the VSIE-diaoptics method (note that in the diaoptics approach, the largest matrix equation that has to be solved is of the size N_s) and $M_{\text{tot}}^{\text{VIE}} = 764$ MB for the MoM-VIE method. To validate the diaoptics solution further, Fig. 3(b) also shows its excellent agreement with the RCS computed by means of the pure MoM-SIE method (with LU decomposition) [7], which takes $N_{\text{tot}}^{\text{SIE}} = 6,144$ unknowns, $T_{\text{tot}}^{\text{SIE}} = 170$ s of computation time, and $M_{\text{tot}}^{\text{SIE}} = 604$ MB of RAM.

B. Array of Homogeneous Dielectric Spherical Scatterers

As an example of curved structures, we replace the cubical scatterers in Fig. 3(a) by spherical ones, as portrayed in Fig. 4(a), where $a = \lambda_0/2.31$ (sphere diameter), $d = \lambda_0/4.29$, $\epsilon_r = 4$, $\theta_{\text{inc}} = 90^\circ$, and $\phi_{\text{inc}} = 0^\circ$. Each diaoptics subsystem, in spite of curvature, is again represented using a single VIE generalized hexahedron with $K_u = K_v = K_w = 2$ and $N_u = N_v = N_w = 4$, bounded by a diaoptics surface constructed from six SIE generalized quadrilaterals with $K_u = K_v = 2$ and $N_u = N_v = 2$, so the numbers of unknowns are $N_v = 240$, $2N_s = 6,144$, and $N_{\text{tot}}^{\text{Dia}} = 6,384$, computation time is $T_{\text{tot}}^{\text{Dia}} = 174$ s, and computer memory consumption is $M_{\text{tot}}^{\text{Dia}} = 152$ MB. The agreement of the RCS results with both the pure MoM-VIE ($N_{\text{tot}}^{\text{VIE}} = 15,360$, $T_{\text{tot}}^{\text{VIE}} = 2,999$ s, $M_{\text{tot}}^{\text{VIE}} = 3.77$ GB) and MoM-SIE solutions is, as can be observed from Fig. 4(b), again excellent, and the computation time with the diaoptics method is reduced by 17.2 times and the memory consumption by 24.8 times when compared to the MoM-VIE method. The MoM-SIE simulation requires

$N_{\text{tot}}^{\text{SIE}} = 6,144$, $T_{\text{tot}}^{\text{SIE}} = 333$ s, and $M_{\text{tot}}^{\text{SIE}} = 604$ MB.

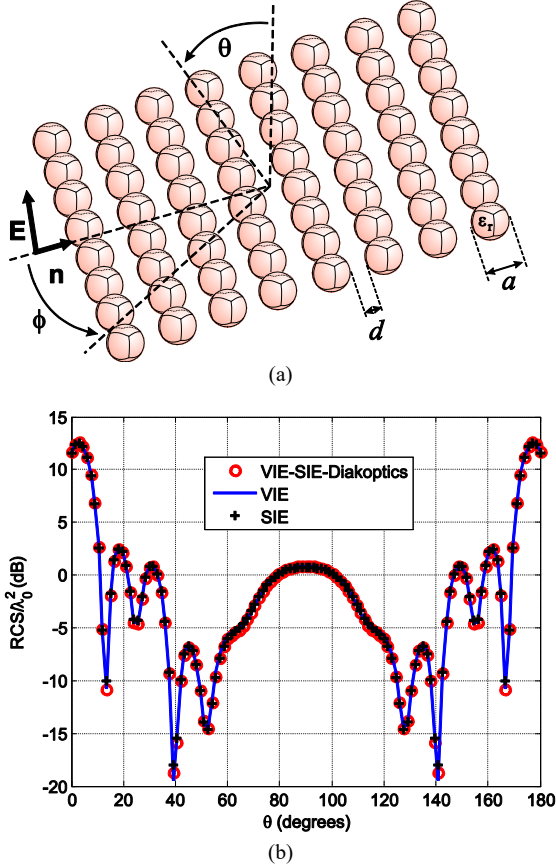


Fig. 4. (a) Array of homogeneous dielectric spherical scatterers excited by a uniform plane wave ($\theta_{\text{inc}} = 90^\circ$, $\phi_{\text{inc}} = 0^\circ$) and (b) normalized bistatic RCS results in the $\phi = 90^\circ$ cut, obtained by the VSIE-diakoptic, pure MoM-VIE, and pure MoM-SIE method.

C. Array of Continuously Inhomogeneous Spherical Scatterers

As an example of inhomogeneous structures, that are also curved, we next consider a 4×4 array of continuously inhomogeneous dielectric spherical scatterers of diameters $a = \lambda_0/1.5$, shown in Fig. 5(a). All sphere-to-sphere distances are $d = \lambda_0/3$ and $\theta_{\text{inc}} = 0^\circ$ and $\phi_{\text{inc}} = 0^\circ$ for the impinging uniform plane wave. The inhomogeneity consists of a linear radial variation of the relative permittivity from $\epsilon_r = 1$ at the surface to $\epsilon_r = 6$ at the center of the spheres, as depicted in the inset of Fig. 5(b). Each scatterer is modeled by seven curvilinear hexahedral VIE elements (with $K_u = K_v = K_w = 2$), one element (with $N_u = N_v = N_w = 2$) approximating the central sphere (of diameter $a/20$) and six continuously inhomogeneous cushion-like elements (with $N_u = N_v = N_w = 3$) attached to the central element across the six corresponding sides, which can be seen in Fig. 5(a).

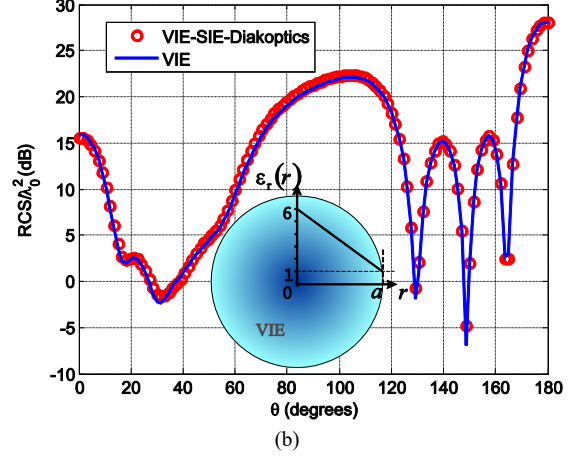
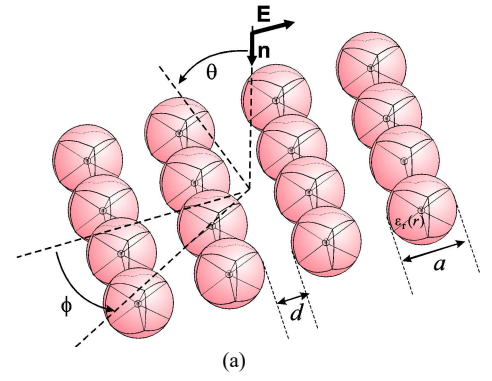


Fig. 5. (a) Array of continuously inhomogeneous dielectric spherical scatterers and (b) bistatic RCS results in the $\phi = 90^\circ$ cut obtained by the VSIE-diakoptic and pure MoM-VIE methods.

The diakoptic surface enclosing each subsystem consists of six SIE patches with $K_u = K_v = 2$ and $N_u = N_v = 4$. The resulting numbers of unknowns are $N_V = 630$ (per subsystem), $2N_S = 6,144$ (total), and $N_{\text{tot}}^{\text{Dia}} = 6,774$, and the total computation time is $T_{\text{tot}}^{\text{Dia}} = 306$ s. In Fig. 5(b), an excellent agreement of the RCS calculated using the diakoptic method and the results obtained by the pure MoM-VIE method, with $N_{\text{tot}}^{\text{VIE}} = 10,080$ and $T_{\text{tot}}^{\text{VIE}} = 23,983$ s, is observed. In this case, the acceleration is even more dramatic (about 78.4 times) than in previous examples, because the MoM-VIE computation of matrix entries for a subsystem involving several inhomogeneous hexahedra is considerably more time consuming than for subsystems with a single homogeneous element. The required RAM for the VSIE-diakoptic and MoM-VIE solutions is $M_{\text{tot}}^{\text{Dia}} = 162$ MB and $M_{\text{tot}}^{\text{VIE}} = 1.6$ GB, respectively.

D. Diakoptic Decomposition of a Large Dielectric Slab

As an example of a single solid object decomposed into diakoptic subsystems, consider scattering from a dielectric slab with relative permittivity $\epsilon_r = 2.25$ and dimensions $6\lambda_d \times 6\lambda_d \times 2\lambda_d$, with λ_d standing for the wavelength in the dielectric. The plane wave excitation is as in the previous

examples ($\theta_{\text{inc}} = 90^\circ$, $\phi_{\text{inc}} = 0^\circ$). The slab is decomposed into nine cubical diakoptic subsystems of edge lengths $d = 2\lambda_d$, as shown in Fig. 6(a). A single VIE element is used for each subsystem, with $K_u = K_v = K_w = 1$ and $N_u = N_v = N_w = 9$ ($N_v = 2,430$ per subsystem); it is enclosed by a six-patch diakoptic surface, with $K_u = K_v = 1$ and $N_u = N_v = 6$ (total of $2N_s = 7,776$ diakoptic unknowns). Note that the common surface between the neighboring diakoptic subsystems is taken into account by finding the limit of SIE integrals for the case when two (diakoptic) surfaces are infinitesimally close.

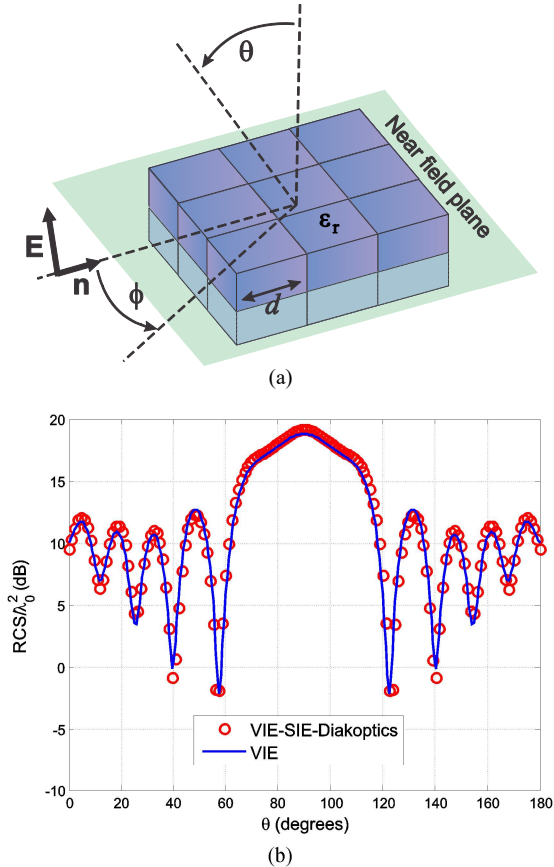


Fig. 6. (a) Large dielectric slab decomposed into nine diakoptic subsystems and (b) comparison of the VSIE-diakoptic solution for the normalized bistatic RCS, in the $\phi = 0^\circ$ cut, of the slab with results obtained using the pure MoM-VIE method.

In Fig. 6(b), we observe an excellent agreement of the VSIE-diakoptic RCS results for the dielectric slab ($N_{\text{tot}}^{\text{Dia}} = 10,206$, $T_{\text{tot}}^{\text{Dia}} = 118$ s, $M_{\text{tot}}^{\text{Dia}} = 373$ MB) with the solution obtained by the pure MoM-VIE method with $N_u = N_v = N_w = 9$ ($N_{\text{tot}}^{\text{VIE}} = 21,708$, $T_{\text{tot}}^{\text{VIE}} = 3,291$ s, $M_{\text{tot}}^{\text{VIE}} = 7.5$ GB). Hence the diakoptic method is 27.9 times faster and 20.2 times less expensive at memory consumption than the pure MoM approach in this case. For additional validation of the proposed diakoptic method, Fig. 7 shows the near (internal or external) total electric field computed in the plane indicated in Fig. 6(a). The VSIE-diakoptic results are

compared with a pure-MoM-SIE solution ($N_{\text{tot}}^{\text{SIE}} = 3,936$, $T_{\text{tot}}^{\text{SIE}} = 68$ s, $M_{\text{tot}}^{\text{SIE}} = 248$ MB), and we observe an excellent agreement. As expected, the SIE solution, with modeling only the external, dielectric-air, surface of the slab in Fig. 6(a) and no unknowns for modeling the internal boundary surfaces between the adopted diakoptic subsystems, as well as no volume unknowns, is more efficient than the VSIE-diakoptic solution in this example. Note, however, that the results in this example and all other examples in this paper are aimed to demonstrate considerably higher efficiency of the VSIE-diakoptic method when compared to the pure MoM-VIE technique implementing the same discretization, with the pure MoM-SIE method serving just for validation and its numerical performance parameters being given for reference.

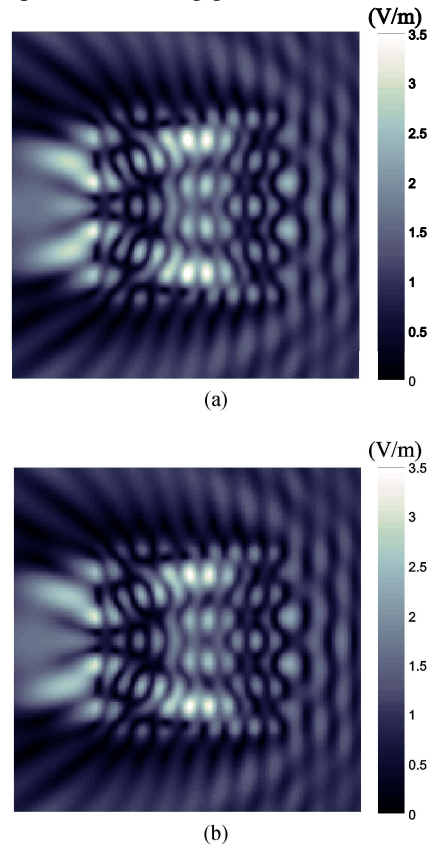


Fig. 7. Magnitude of the near total electric field inside and around the dielectric slab scatterer computed in the plane indicated in Fig. 6(a), where the size of the near-field computation area is $11\lambda_d \times 11\lambda_d$, and it cuts through the middle of the vertical dimension of the slab: comparison of (a) VSIE-diakoptic results and (b) pure MoM-SIE reference solution.

E. Diakoptic Decomposition of a Human Phantom

Finally, as an example of a complex object decomposed into completely different subsystems, consider scattering from a human phantom. The excitation of the phantom is by a uniform plane wave incident along the x -direction ($\theta_{\text{inc}} = 90^\circ$, $\phi_{\text{inc}} = 0^\circ$) at a frequency of 900 MHz, and the averaged equivalent complex relative permittivity of the homogeneous lossy dielectric material filling the phantom at this frequency is taken to be $\epsilon_{\text{rc}} = 52.72 - j19.66$. The phantom is

decomposed into four diakoaptic subsystems as shown in the inset of Fig. 8. Subsystems are different (the phantom is not symmetric) and are modeled by 1436, 1310, 1050, and 985 VIE hexahedra and 1124, 1114, 1018, and 976 SIE quadrilaterals, respectively (a total of $2N_S = 16,928$ diakoaptic unknowns). Initial triangular mesh, provided by NEVA Electromagnetics [42], is re-meshed in ANSYS ICEM CFD 15.0.

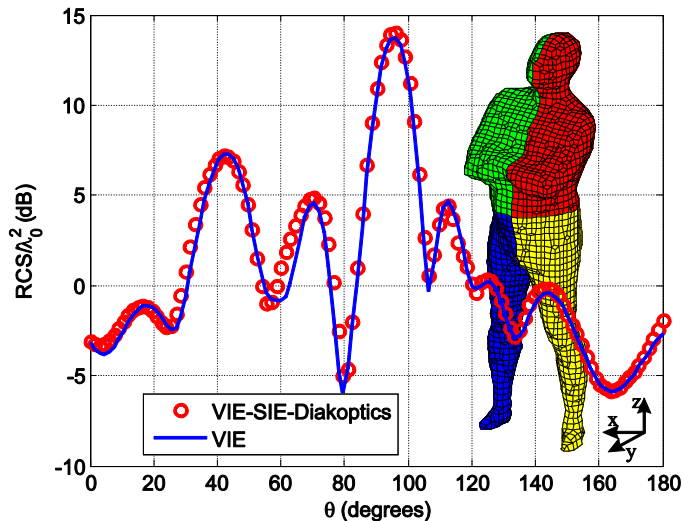


Fig. 8. Analysis of a human phantom decomposed into four diakoaptic subsystems (figure inset): comparison of the VSIE-diaokoaptic solution for the normalized bistatic RCS ($\phi = 0^\circ$ cut) of the phantom with results obtained using the pure MoM-VIE method.

In Fig. 8, a good agreement of the VSIE-diaokoaptic RCS results for the human phantom ($N_{\text{tot}}^{\text{Dia}} = 33,387$, $T_{\text{tot}}^{\text{Dia}} = 17,584$ s, $M_{\text{tot}}^{\text{Dia}} = 3.7$ GB) with the pure MoM-VIE solution ($N_{\text{tot}}^{\text{VIE}} = 16,012$, $T_{\text{tot}}^{\text{VIE}} = 135,256$ s, $M_{\text{tot}}^{\text{VIE}} = 4.1$ GB) is observed. Hence, the acceleration gained with diakoaptics in this example is 7.7 times, and because of the requirement for storage of all four different subsystem matrices, the reduction in RAM consumption is only 11%. Note that the number of VIE unknowns for the full system ($N_{\text{tot}}^{\text{VIE}}$) is smaller than the number of diakoaptic unknowns, $2N_S = 16,928$. In this example, there are two advantages of the VSIE-diaokoaptics approach. On one hand, in diakoaptics, the largest matrix that has to be LU decomposed is of size $N_S = 8,464$ (see Appendix I). On the other, the burdensome computation of the interaction of VIE elements belonging to different subsystems is replaced with faster computation of interaction between diakoaptic surfaces.

IV. CONCLUSIONS

This paper has proposed a novel diakoaptic method for 3-D electromagnetic analysis based on MoM-VSIE modeling. The method breaks the original structure into a number of non-overlapping closed-region subsystems (which may be the same or different) containing inhomogeneous dielectric

materials and an open-region subsystem that encloses all other subsystems. Each subsystem is analyzed completely independently applying the double-higher-order large-domain Galerkin generalized MoM-VSIE or MoM-SIE solvers, and the final solution is obtained from diakoaptic matrices expressing linear relations between electric and magnetic equivalent surface current coefficients on diakoaptic surfaces.

The proposed VSIE-diaokoaptic method has been demonstrated, evaluated, and discussed in several characteristic examples, including both homogeneous and continuously inhomogeneous dielectric structures, objects with both flat surfaces/sharp edges and pronounced curvature, and both far-field and near-field computations. Numerical examples have demonstrated that the proposed VSIE-diaokoaptic method substantially increases the efficiency of the conventional pure MoM-VIE approach, with the pure MoM-SIE method serving for validation, as a very different technique. When compared to the pure MoM-VIE technique implementing the same type of discretization, the diakoaptic approach enables considerable accelerations and memory savings, while fully preserving the accuracy of the analysis. The paper has also demonstrated that the general diakoaptic methodology can effectively include VIE-based subsystems and has shown how that can be done.

APPENDIX I

A. Theoretical Analysis of Computational Complexity and Acceleration of VSIE-Diaokoaptics

Here, we make a theoretical estimate for the maximum efficiency of the diakoaptic analysis, as compared to the standard MoM-VIE analysis, assuming single-level decomposition. We also assume that it is possible to subdivide the EM system into M congruent (i.e., geometrically identical) diakoaptic subsystems, which differ only in their (physical) position in the original system and are analyzed by the generalized VSIE method. These subsystems are “wrapped” with a total of MN_D unknowns, namely, each subsystem has $N_D = 2N_S$ SIE unknowns on its surface, with N_S standing for the number of unknown coefficients for expansion of each of the equivalent currents \mathbf{J}_e and \mathbf{M}_e on the surface. Every VIE subsystem has N_V volumetric unknowns for the approximation of the displacement vector \mathbf{D} , so the total of VIE unknowns is MN_V . We assume, further, classical MoM codes, which utilize a direct solver based on LU factorization, so that the matrix filling and matrix solution have computational time complexities of $O(N^2)$ and $O(N^3)$, respectively, where N is the size of the matrix. Hence, the total analysis time of the entire VIE problem solved at once is given by $T_{\text{tot}}^{\text{VIE}} = k_V(MN_V)^2 + k^{\text{solV}}(MN_V)^3$, with k_V and k^{solV} standing for averaged times spent for each procedure per unknown coefficient. The total VIE filling and solution times are $k_V N_V^2$ and $k^{\text{solV}} N_V^2$, and those for the SIE are $k_S N_D^2$ and $k^{\text{solV}} N_D^2$, respectively. Therefore, the filling time for each

VSIE subsystem matrix can be computed as $k_V^{\text{sub}} N_V^2 + k_S^{\text{sub}} N_D^2 + 2k_{VS}^{\text{sub}} N_V N_D$, where k_S^{sub} , k_{VS}^{sub} , and k_V^{sub} are averaged times spent per computation of the corresponding matrix element in the subsystem.

In the proposed VSIE-diaoptical approach, matrices to be computed are of sizes $N_D + N_V$ and MN_D , and the sizes of matrices that need to be solved are MN_S and $N_V + N_S$. Taking into account that $k_S^{\text{sub}} \ll k_{VS}^{\text{sub}} \ll k_V^{\text{sub}}$ (since k_S^{sub} , k_{VS}^{sub} , and k_V^{sub} are essentially the times required for surface-surface, surface-volume, and volume-volume integration, respectively), the total time needed for the diaoptical analysis is

$$T_{\text{tot}}^{\text{Dia}} \approx \max \left\{ k_V^{\text{sub}} N_V^2, k_S \left(2MN_S \right)^2 \right\} + k^{\text{solv}} \left[2 \left(MN_S \right)^3 + \left(N_V + N_S \right)^3 \right] \quad (15)$$

Note that $k_V^{\text{sub}} \neq k_V$, which is because the computation of matrix elements obtained via integration of basis functions belonging to the same element or neighboring elements requires more time than for those related to the distant elements. In our numerical code, the average filling time of the entire VIE system is related to the time for its subsystem as $k_V^{\text{sub}} = k_V \sqrt[3]{M}$.

Based on the above, the theoretically maximum acceleration of the VSIE-diaoptics as compared to the standard MoM-VIE analysis can be estimated as

$$\frac{T_{\text{tot}}^{\text{VIE}}}{T_{\text{tot}}^{\text{Dia}}} \approx \frac{k_V \left(MN_V \right)^2 + k^{\text{solv}} \left(MN_V \right)^3}{T_{\text{fill}}^{\text{Dia}} + k^{\text{solv}} \left[2 \left(MN_S \right)^3 + \left(N_V + N_S \right)^3 \right]}, \quad (16)$$

where $T_{\text{fill}}^{\text{Dia}} = \max \{ k_V \sqrt[3]{M} N_V^2, k_S \left(2MN_S \right)^2 \}$ is the time required for filling of all diaoptical matrices, computed simultaneously. Additionally, (16) can be generalized to predict acceleration of the system subdivided into non-congruent subsystems. In such a case, $T_{\text{fill}}^{\text{Dia}}$ is the maximum of all anticipated filling times for the diaoptical matrices. Table I gives the actually measured and theoretically maximum acceleration values for all examples in the paper, in Sections III.A–E.

Note that the anticipated theoretical acceleration in (16) is the maximum acceleration that can possibly be achieved, by the respective system subdivision using a given numerical VSIE-diaoptical code. The actual acceleration is lower, due to the fact that the computation time $T_{\text{tot}}^{\text{Dia}}$ does not include times required for preprocessing, post-processing, and matrix assembly (summation of diaoptical subsystem sub-matrices and row/column manipulations), as well as the

$k_S^{\text{sub}} N_D^2 + 2k_{VS}^{\text{sub}} N_V N_D$ portion of the subsystem matrix filling time. However, despite of all these omissions and simplifications, the simple formula in (16) provides, as can be observed from Table I, a good estimate of the efficiency of the diaoptical performance. Namely, all actually measured acceleration values agree well with the corresponding theoretically maximum values in Table I – for all examples in the paper.

We also observe from Table I considerable (measured and theoretical) accelerations, while fully preserving the accuracy of the analysis, enabled by the diaoptical approach when compared to the pure MoM-VIE technique implementing the same type of discretization, in all examples in Sections III.A–E.

TABLE I
ACTUALLY MEASURED AND THEORETICALLY MAXIMUM ACCELERATION

Example in Sections III.A–E	Measured Acceleration	Theoretically Max. Acceleration
Array of Homogeneous Dielectric Cubical Scatterers	4.5	5.9
Array of Homogeneous Dielectric Spherical Scatterers	17.2	20.1
Array of Continuously Inhomogeneous Spherical Scatterers	78.4	95.1
Diaoptical Decomposition of a Large Dielectric Slab	27.9	31.3
Diaoptical Decomposition of a Human Phantom	7.7	9.4

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