

# MoM/BI-RME Analysis of Boxed MMICs With Arbitrarily Shaped Metallizations

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**Abstract**—In this paper, we propose a novel approach for the analysis of shielded microstrip circuits, composed of a number of thin metallic areas with arbitrary shapes and finite conductivity, embedded in a multilayered lossy medium. The analysis is based on the solution of an integral equation (IE) obtained by enforcing the proper boundary condition to the electric field on the metallic areas. The IE is solved by using the method of moments with entire domain basis functions, which are numerically determined by the boundary integral-resonant-mode expansion (BI-RME) method. The use of the BI-RME method allows for the efficient calculation of the basis functions independently on the shape of the domain, thus permitting the analysis of a wide class of circuits. Two examples demonstrate the accuracy, rapidity, and flexibility of the proposed method.

**Index Terms**—Entire domain basis functions, integral equations, microstrip filters, MMICs, moment method.

## I. INTRODUCTION

OVER THE last years, considerable interest has been directed to the design of boxed multilayered circuits (Fig. 1). This configuration is typically considered in the design of many actual monolithic microwave integrated circuits (MMICs), both in single-layer [1] and multilayered configurations [2], [3].

Among the possible numerical methods applied to the analysis of this type of structures, the integral equation (IE) method is by far one of the most efficient. The IE method can be formulated either in the spectral [4] or spatial domains [5]. The resulting IE is solved by the method of moments (MoM), usually considering sub-domain basis functions (e.g., rooftops [6], [7] or basis functions on triangular domains [8], [9]).

Recently, the IE/MoM method in the spectral domain was applied with entire domain basis functions [10]. The main advantage of using the vector modal functions derived in [10] is the dramatic reduction in the order of the MoM matrix since few entire domain basis functions are usually sufficient to represent

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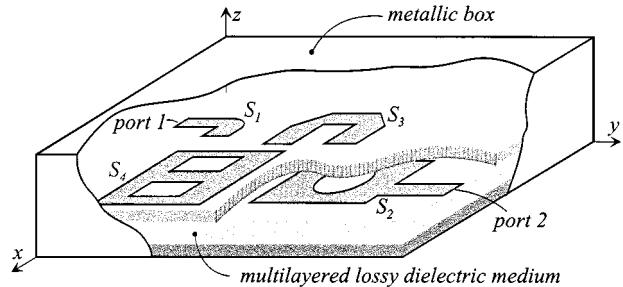


Fig. 1. Shielded MMIC with arbitrarily shaped metallic areas in a multilayered lossy medium. The patches can be placed at different height and they may overlap.

the unknown currents. Moreover, the calculation of the MoM matrix is enhanced because of the rapid decrease of the spectral components to be added subsequently for the calculation of the matrix elements. It is noted, however, that the original work derived in [10] is limited to areas with a rectangular shape, where the entire domain basis functions are known analytically.

In this paper, we present the extension of the method proposed in [10] to the case of metallic areas with an arbitrary shape (Fig. 1). The entire domain basis functions are determined numerically by the boundary integral-resonant mode expansion (BI-RME) method [11]. The use of the BI-RME method has two main advantages. The first is the possibility of obtaining entire domain basis functions for arbitrary shapes in a short time, and the second is that the entries of the MoM matrix are practically obtained as a by-product of the method itself. In fact, the surface integrals involved in the calculations of the MoM matrix can be converted into line integrals on the boundary of the metallic areas, and the quantities required on the boundary are the basic output of the BI-RME calculation.

A preliminary discussion of the proposed algorithm was presented in [12]. This paper gives a comprehensive explanation of the MoM/BI-RME method, adding two novel capabilities: metallic areas including a port may have an arbitrary shape, and multiply connected metallizations can be considered.

## II. IE/MoM APPROACH

Let us consider the structure shown in Fig. 2, consisting of a multilayered medium and  $P$  metallic areas with arbitrary shapes  $S_1, \dots, S_P$ , possibly located at different interfaces. The circuit is fed at the frequency  $\omega$  at  $K$  ports ( $K \leq P$ ), conventionally defined on the first  $K$  areas  $S_1, \dots, S_K$ . As usual [5], [13], [14], the ports are represented as small gaps between the metallization and the wall of the box (delta-gap voltage excitations).

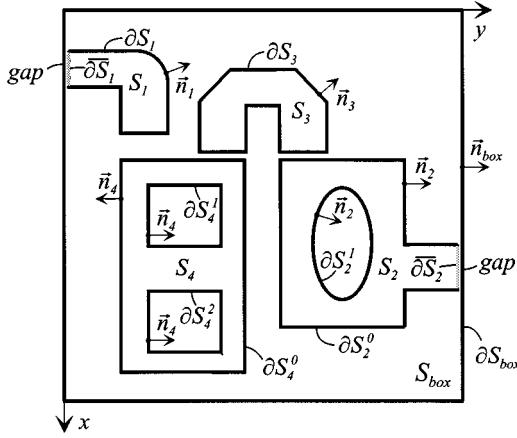


Fig. 2. Different metallic areas considered in the analysis: simply ( $S_1$ ) or multiply connected patches attached to a port ( $S_2$ ), simply ( $S_3$ ) or multiply connected internal patches ( $S_4$ ).  $\partial S_p$  represents the boundary of the area  $S_p$ , except the possible part corresponding to a port (denoted by  $\overline{\partial S}_p$ ). In case of multiply connected areas,  $\partial S_p$  consists of many lines  $\partial S_p^0$ ,  $\partial S_p^1$ , ...

In [10], the analysis is based on the solution of a system of  $P$  IEs, which are obtained by enforcing the boundary condition to the transverse-to- $z$  electric field at all the metallic areas

$$Z \vec{J}_p(\vec{r}) - \sum_{q=1}^P \int_{S_q} \vec{G}(\vec{r}, \vec{r}'|\omega) \cdot \vec{J}_q(\vec{r}') dS' = \begin{cases} -\vec{n}_p v_p \delta(\vec{r}, \vec{r}_p)|_{\vec{r}_p \in \overline{\partial S}_p}, & p = 1, \dots, K \\ 0, & p = K+1, \dots, P \end{cases} \quad (1)$$

where the points  $\vec{r}$  and  $\vec{r}'$  belong to  $S_p$  and  $S_q$ , respectively,  $\overline{\partial S}_p$  denotes the part of the boundary of  $S_p$  that belongs to the  $p$ th port,  $\vec{n}_p$  is the outward normal,  $v_p$  is the voltage applied to the  $p$ th port, and  $Z$  is the “sheet impedance” of the metallizations. The choice of  $Z$  is problem dependent. For instance, in the case of a single-layer microstrip circuit with a metallization thickness  $t$  much larger than the skin depth  $\delta$ , we can use the surface impedance of the conductors  $Z = Z_s = (1+j)\rho/\delta$  ( $\rho$  is the resistivity of the metal), in the case of low-frequency calculations ( $t \ll \delta$ ), we can use  $Z = \rho/t$ . Moreover,  $\vec{J}_q$  is the (unknown) current density on  $S_q$ , and the Green's function  $\vec{G}$  is given by [15]

$$\vec{G}(\vec{r}, \vec{r}'|\omega) = \sum_m V_m(z, z'|\omega) \vec{\mathcal{E}}_m(x, y) \vec{\mathcal{E}}_m(x', y') \quad (2)$$

where  $\vec{\mathcal{E}}_m$  are the transverse electric modal vectors of the TE and TM modes of the box, with the normalization  $\int_{S_{box}} |\vec{\mathcal{E}}_m|^2 dS = 1$ . The expressions of  $\vec{\mathcal{E}}_m$  are

$$\vec{\mathcal{E}}_m' = -\frac{\nabla_T \chi_m'}{k_m'} \quad (\text{TM modes}) \quad (3)$$

$$\vec{\mathcal{E}}_m'' = -\vec{u}_z \times \frac{\nabla_T \chi_m''}{k_m''} \quad (\text{TE modes}) \quad (4)$$

where  $\chi_m'$  and  $\chi_m''$  are the eigenfunctions of the Helmholtz equation with the Dirichlet or Neumann boundary conditions on

$\partial S_{box}$  (see Fig. 2), and  $k_m'$  and  $k_m''$  are the corresponding eigenvectors. Finally, functions  $V_m$  are determined by considering the equivalent modal transmission lines for the layered box [10], [16].

Equation (1) is solved by applying the MoM in the Galerkin form. The unknown current density  $\vec{J}_p$  is represented through a suitable set of  $N_p$  basis functions  $\vec{e}_r^{(p)}$  defined on the  $p$ th patch, namely,

$$\vec{J}_p = \sum_{r=1}^{N_p} \xi_r^{(p)} \vec{e}_r^{(p)}, \quad p = 1, \dots, P \quad (5)$$

where  $\xi_r^{(p)}$  are unknown coefficients.

As discussed in [10], the calculation of the MoM matrices involves frequency-independent coefficients of two types. The former represents the *coupling integral* between the  $r$ th basis function on the  $p$ th metallic area and the  $m$ th modal vector of the box, and is given by

$$\mathcal{C}_{rm}^{(p)} = \int_{S_p} \vec{e}_r^{(p)}(\vec{r}) \cdot \vec{\mathcal{E}}_m(\vec{r}) dS. \quad (6)$$

The latter is the projection of the delta-gap excitation of the  $p$ th port on the  $r$ th basis function (*port integral*), and is given by

$$\mathcal{P}_r^{(p)} = \int_{\overline{\partial S}_p} \vec{e}_r^{(p)}(\vec{r}) \cdot (-\vec{n}_p) d\ell. \quad (7)$$

For any frequency, the scattering parameters are calculated straightforwardly from the coefficients  $\xi_r^{(p)}$  obtained by the solution of the MoM system [10].

### III. ENTIRE DOMAIN BASIS FUNCTIONS

A key feature of the present approach is the use of a set of entire domain basis functions, i.e., functions  $\vec{e}_r^{(p)}$ , which span the entire domain  $S_p$ .

The advantage of using such functions has been demonstrated in [10], with reference to the case where all surfaces  $S_p$  are rectangular. More specifically, the electric modal vectors of rectangular waveguides bounded by magnetic or mixed-type walls have been used as basis functions. In this paper, the same concept is applied to metallization of arbitrary shapes, with no restriction on the geometry of the surfaces  $S_p$ . In this case, the basis functions must be determined numerically and the efficiency of the numerical method used for their calculation is a vital issue. The BI-RME method discussed in Section V permits to determine very efficiently enough basis functions for our application.

In the case of simply connected surfaces, we have to calculate two classes of basis functions expressed by

$$\vec{e}_r'^{(p)} = -\vec{u}_z \times \frac{\nabla_T \psi_r'^{(p)}}{\kappa_r'} \quad (8)$$

$$\vec{e}_r''^{(p)} = -\frac{\nabla_T \psi_r''^{(p)}}{\kappa_r''^{(p)}} \quad (9)$$

where the pairs  $\{\psi_r'^{(p)}, \kappa_r'^{(p)}\}$  and  $\{\psi_r''^{(p)}, \kappa_r''^{(p)}\}$  are the eigen-solutions of the homogeneous Helmholtz equation in the domain  $S_p$ , i.e.,

$$\nabla_T^2 \psi_r'^{(p)} + \kappa_r'^{(p)2} \psi_r'^{(p)} = 0, \quad \text{in } S_p \quad (10)$$

$$\nabla_T^2 \psi_r''^{(p)} + \kappa_r''^{(p)2} \psi_r''^{(p)} = 0, \quad \text{in } S_p. \quad (11)$$

In the case of  $N$ -times connected surfaces, the set of basis functions must be supplemented with  $N - 1$  additional functions

$$\vec{e}_r^{0(p)} = -\vec{u}_z \times \nabla_T \psi_r^{0(p)} \quad (12)$$

where  $\psi_r^{0(p)}$  satisfies the Laplace equation in the domain  $S_p$ , i.e.,

$$\nabla_T^2 \psi_r^{0(p)} = 0, \quad \text{in } S_p. \quad (13)$$

The boundary conditions are different for metallizations connected or not connected to ports.

In the case of metallizations not connected to ports (e.g.,  $S_3$  and  $S_4$  in Fig. 2),  $\vec{J}_p$  is tangent to the whole boundary of  $S_p$ . The basis functions  $\vec{e}_r'^{(p)}$ ,  $\vec{e}_r''^{(p)}$ , and  $\vec{e}_r^{0(p)}$  satisfy the same boundary condition provided that

$$\psi_r'^{(p)} = 0, \quad \text{on } \partial S_p \quad (14)$$

$$\partial \psi_r''^{(p)} / \partial n_p = 0, \quad \text{on } \partial S_p \quad (15)$$

$$\psi_r^{0(p)} = \begin{cases} 1, & \text{on the inner contour } \partial S_p^i \\ 0, & \text{on } \partial S_p - \partial S_p^i \end{cases} \quad (16)$$

In the case of metallizations connected to ports (e.g.,  $S_1$  and  $S_2$  in Fig. 2),  $\vec{J}_q$  is perpendicular to the port segment  $\overline{\partial S_p}$ . In this case,  $\psi_r'^{(p)}$ ,  $\psi_r''^{(p)}$ , and  $\psi_r^{0(p)}$  must satisfy the mixed boundary conditions

$$\begin{cases} \psi_r'^{(p)} = 0, & \text{on } \partial S_p \\ \partial \psi_r'^{(p)} / \partial n_p = 0, & \text{on } \overline{\partial S_p} \end{cases} \quad (17)$$

$$\begin{cases} \partial \psi_r''^{(p)} / \partial n_p = 0, & \text{on } \partial S_p \\ \psi_r''^{(p)} = 0, & \text{on } \overline{\partial S_p} \end{cases} \quad (18)$$

$$\begin{cases} \psi_r^{0(p)} = \begin{cases} 1, & \text{on the inner contour } \partial S_p^i \\ 0, & \text{on } \partial S_p - \partial S_p^i \end{cases} \\ \partial \psi_r^{0(p)} / \partial n_p = 0, & \text{on } \overline{\partial S_p} \end{cases} \quad (19)$$

It is worthy noting that, in the case of metallizations connected to ports,  $\partial S_p$  denotes the boundary of  $S_p$ , but the port segment  $\overline{\partial S_p}$ .

#### IV. COUPLING AND PORT INTEGRALS

When considering metallic areas with a rectangular shape, the coupling integrals (6) and the port integrals (7) can be calculated analytically by using the analytical expressions of the basis functions (see, for instance, [10]). This possibility is precluded in the case of arbitrary shapes because the basis functions are determined numerically. In this case, the surface integration (6) is a time-consuming task, especially in cases of basis functions determined by a boundary integral method, since it requires the numerical evaluation of the basis functions in many points within the integration domain  $S_p$ . However, the coupling integrals (6) can be transformed from surface to line integrals,

thus dramatically reducing the computing time. As shown in the Appendix, we have

$$\int_{S_p} \vec{e}_r'^{(p)} \cdot \vec{\mathcal{E}}_m' dS = 0 \quad (20)$$

$$\int_{S_p} \vec{e}_r''^{(p)} \cdot \vec{\mathcal{E}}_m' dS = \frac{\kappa_r''^{(p)}}{k_m' \left( \kappa_r''^{(p)2} - k_m'^2 \right)} \int_{\partial S_p} \psi_r''^{(p)} \frac{\partial \chi_m'}{\partial n_p} d\ell \quad (21)$$

$$\int_{S_p} \vec{e}_r^{0(p)} \cdot \vec{\mathcal{E}}_m' dS = 0 \quad (22)$$

$$\int_{S_p} \vec{e}_r'^{(p)} \cdot \vec{\mathcal{E}}_m'' dS = \frac{k_m''}{\kappa_r'^{(p)} \left( \kappa_r'^{(p)2} - k_m''^2 \right)} \int_{\partial S_p} \chi_m'' \frac{\partial \psi_r'^{(p)}}{\partial n_p} d\ell \quad (23)$$

$$\int_{S_p} \vec{e}_r''^{(p)} \cdot \vec{\mathcal{E}}_m'' dS = \frac{1}{\kappa_r''^{(p)} k_m''} \int_{\partial S_p} \psi_r''^{(p)} \frac{\partial \chi_m''}{\partial t_p} d\ell \quad (24)$$

$$\int_{S_p} \vec{e}_r^{0(p)} \cdot \vec{\mathcal{E}}_m'' dS = -\frac{1}{k_m''} \int_{\partial S_p} \chi_m'' \frac{\partial \psi_r^{0(p)}}{\partial n_p} d\ell \quad (25)$$

where  $\partial / \partial t_p$  is the derivative along the boundary, namely, in the direction of  $\vec{t}_p = \vec{u}_z \times \vec{n}_p$ .

It is noted that these formulas hold true in both cases of metallizations connected or not connected to ports.

For the port integrals (7), we easily obtain

$$\int_{\overline{\partial S_p}} \vec{e}_r'^{(p)}(\vec{r}) \cdot (-\vec{n}_p) d\ell = 0 \quad (26)$$

$$\int_{\overline{\partial S_p}} \vec{e}_r''^{(p)}(\vec{r}) \cdot (-\vec{n}_p) d\ell = \frac{1}{\kappa_r''^{(p)}} \int_{\overline{\partial S_p}} \frac{\partial \psi_r''^{(p)}}{\partial n_p} d\ell \quad (27)$$

$$\int_{\overline{\partial S_p}} \vec{e}_r^{0(p)}(\vec{r}) \cdot (-\vec{n}_p) d\ell = 0. \quad (28)$$

#### V. APPLICATION OF THE BI-RME METHOD

In the analysis of circuits of practical interest, some tens of basis functions (8) and (9) are usually needed for each metallic area. This, in turn, requires the calculation of some tens of eigen-solutions of Helmholtz equations (10) and (11).

In the past years, some of the authors developed a novel method (i.e., the BI-RME method) for the solution of the Helmholtz equation in arbitrary domains [17]–[19]. A comprehensive description of the BI-RME method is reported in [11]. In this section, we limit ourselves to a brief outline of the BI-RME method, to highlight its advantages in the calculation of entire domain basis functions.

The BI-RME method is a modified boundary integral approach for the evaluation of eigenfunctions. The surface  $S_p$  (either simply or multiply connected) is considered as a part of a fictitious enlarged domain  $\Omega_p$  with a rectangular shape [see Fig. 3(a)]. The eigenfunctions to be determined are defined in  $\Omega_p$ , and are assumed to vanish outside  $S_p$ . They are expressed as combinations of boundary integrals (BIs) and a resonant mode expansion (RME), involving the modal potentials of the region  $\Omega_p$ . The boundary integrals involve  $\partial \psi_r'^{(p)} / \partial n_p$  and  $\psi_r''^{(p)}$  over the line  $\sigma_p$  [see Fig. 3(a)], which corresponds to the part of  $\partial S_p$  not coincident with the rectangular boundary. Using

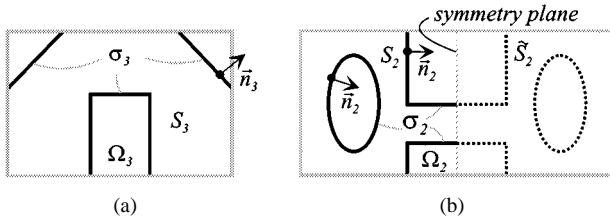


Fig. 3. Geometry for the application of the BI-RME method. (a) Internal metallic area. (b) Metallic area connected to a port.

the BI-RME representation of the eigenfunctions and imposing the proper boundary conditions on  $\sigma_p$ , the eigenvalue problems (10) and (11) are converted into integral-differential equations. As in the case of the conventional boundary element method (BEM), the discretized problem resulting from the application of the BI-RME method is much smaller than in conventional approaches based on differential equations (finite element method (FEM), finite-difference time-domain (FDFD), etc.). Differently from the conventional BEM, the BI-RME method leads to the determination of the eigenfunctions by the solution of a linear matrix eigenvalue problem. In particular, it provides as eigenvalues  $\kappa_r^{(p)}$  and  $\kappa_r^{\prime\prime(p)}$  up to a prescribed value  $\kappa_{\max}$ , and as eigenfunctions  $\partial\psi_r^{(p)}/\partial n_p$  and  $\psi_r^{\prime\prime(p)}$  over the line  $\sigma_p$  and the modal amplitudes of the RME.

The method is very efficient and reliable, also in cases where a large number of eigenfunctions have to be determined. Moreover, no spurious modes are found. Furthermore, it is worth noting that the order of the matrix eigenvalue problem to be solved depends on the extension of the line  $\sigma_p$  and of the surface of the resonator  $\Omega_p$ . For this reason, the efficiency of the BI-RME method highly improves when a large part of the boundary of  $S_p$  fits with the rectangular boundary, as in the example in Fig. 3(a).

The evaluation of coupling integrals (21), (23), and (24) requires  $\partial\psi_r^{(p)}/\partial n_p$  and  $\psi_r^{\prime\prime(p)}$  on  $\partial S_p$ . On the portion  $\sigma_p$ , these quantities are directly the solutions of the BI-RME method. On the other part of  $\partial S_p$ , they are obtained from the BI-RME representation of the eigenfunctions (see [11, eqs. (5.44) and (5.94)]). Even if a post-processing is required for obtaining the boundary values on  $\partial S_p - \sigma_p$ , it is convenient to let  $\partial S_p$  coincide with  $\partial\Omega_p$  as much as possible. In fact, this reduces the dimensions of the eigenvalue problems to be solved, thus increasing the rapidity and accuracy of the solution.

To solve the Helmholtz equation in the case of a metallization connected to a port, the exterior domain  $\Omega_p$  includes not only  $S_p$ , but also its mirror image  $\tilde{S}_p$ , as shown in Fig. 3(b). The problem is solved by imposing an even or an odd symmetry condition with respect to the symmetry plane shown in Fig. 3(b). The details on the implementation of the BI-RME method taking into account the symmetries are discussed in [11, Sec. 5.2.3].

Finally, for multiply connected surfaces, all the possible basis functions (12) must be considered, requiring the determination of all the solutions  $\psi_r^{0(p)}$  of the Laplace equation (13). These basis functions are obtained by solving (13) through the conventional BEM [17]. Even in this case, when the patch is connected to a port, it is possible to solve the Laplace equation by

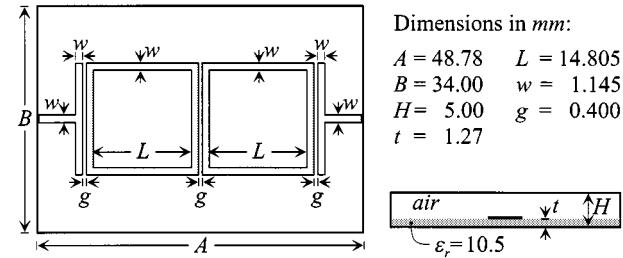


Fig. 4. Printed microstrip filter composed of T-shaped port elements and square-loop resonators.

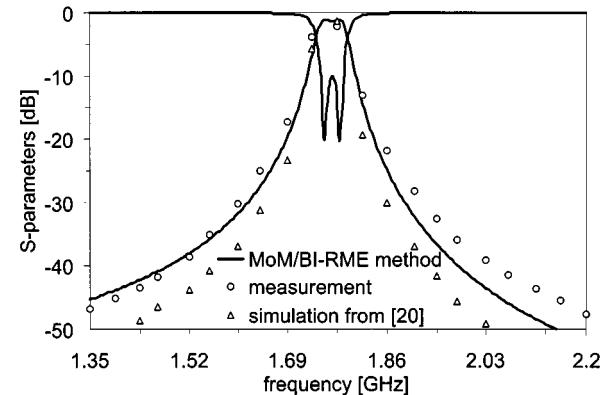


Fig. 5. Comparison between simulated and measured results presented in [20], and results obtained with the present approach, for the filter shown in Fig. 4.

creating a mirror image  $\tilde{S}_p$  of  $S_p$ , and considering an even symmetry condition on the symmetry plane.

## VI. NUMERICAL RESULTS

We used the code for the analysis of printed circuits involving resonators with complex shapes, which fully exploit the capabilities of the method.

The first example refers to the analysis of a narrow-band microstrip filter composed of two T-shaped port elements and two square-loop resonators (Fig. 4), firstly proposed in [20]. The results obtained by the MoM/BI-RME approach are reported in Fig. 5 and compared with experimental data and simulations given in [20], showing a good agreement. The convergence was obtained with 4000 modes of the box and 62 basis functions (eight on each T-shaped line and 23 on each loop resonator), corresponding to  $\kappa_{\max} = 1.153 \text{ mm}^{-1}$ .

For this example, we report in Fig. 6 the study of the convergence properties of the MoM/BI-RME method. In particular, we verified the convergence when varying the number of basis functions [see Fig. 6(a)], and when varying the number of modes of the box [see Fig. 6(b)]. These graphs show that the frequency response does not change when considering more than 62 basis functions or more than 4000 modes of the box.

The calculation of the frequency response of Fig. 5 takes 22 s for the determination of the basis functions and the evaluation of the coupling and port integrals, and 0.135 s for the MoM solution in each frequency point (on a Pentium III at 1 GHz). Thus, the total computing time for the analysis in 100 frequency points is 35.5 s. It is worth observing that taking advantage of the symmetries of the geometry (which were not exploited in

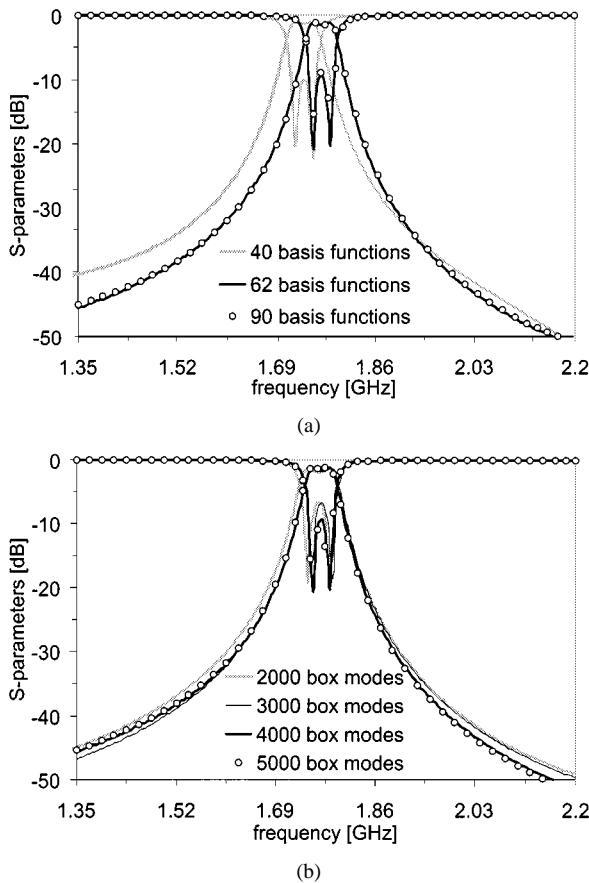


Fig. 6. Convergence behavior of the MoM/BI-RME method in the analysis of the circuit shown in Fig. 4. (a) As a function of the total number of basis functions, when considering 4000 modes of the box. (b) As a function of the number of modes of the box, when considering 62 basis functions.

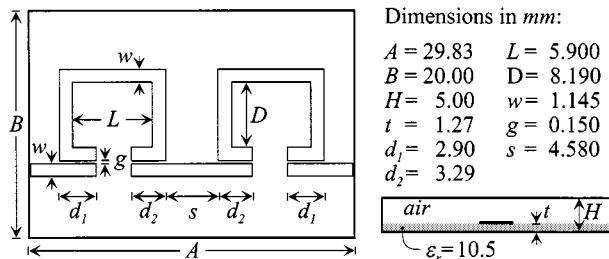


Fig. 7. Printed microstrip filter composed of two open-loop resonators.

our analysis) should lead to a dramatic reduction of the total computing time.

The second example refers to the analysis of a narrow-band microstrip filter composed of two open-loop resonators (Fig. 7) [20]. The analysis of this structure (already presented in [12]) is particularly challenging since the lines and loops are separated by very narrow capacitive gaps, and the surface current must be accurately represented near these gaps. Consequently, in this case, the analysis requires more basis functions than in the previous example. Fig. 8 shows the frequency response of the filter considering 4000 modes of the box and 158 basis functions (12 on each port line, 22 on the middle line, and 56 on each open loop) corresponding to  $\kappa_{\max} = 3.144 \text{ mm}^{-1}$ . The results are in a good agreement with the theoretical and experimental data taken from [20]. The computing time (without exploiting the

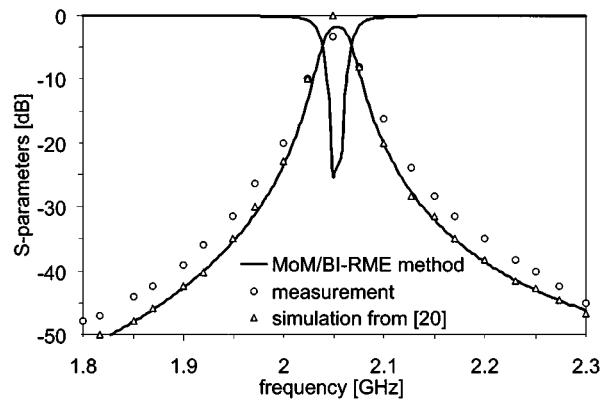


Fig. 8. Comparison between simulated and measured results presented in [20], and results obtained with the present approach, for the filter shown in Fig. 7.

symmetries) is 48 s for the calculation of the basis functions and of the coupling and port integrals, and 0.63 s for the MoM solution in each frequency point (on a Pentium III at 1 GHz). Thus, the total computing time for the analysis in 100 frequency points is 111 s.

As a final remark, we can observe that the selection of the basis functions in our approach is performed by a spectral criterion, including all the entire domain basis functions up to a prescribed  $\kappa_{\max}$ . Especially in cases of metallizations where one of the dimensions is much larger than the other (e.g., narrow strips), this criterion leads to the adoption of very large values of  $\kappa_{\max}$ , in order to include in the set of basis functions a sufficient number of elements with significant variation along the narrow dimension. Of course, it can happen that many basis functions with uselessly rapid variation in the larger direction can be included in the basis. These functions could be discarded, but the procedure for their recognition is too complicate to be conveniently implemented in a general purpose computer code. Anyway, the numerical experiments presented above permit not to dramatize the problem because we noted that it is sufficient to consider basis functions whose variation in the narrow dimension correspond to one or two sinusoidal oscillations. In spite of the roughness of the current representation, the results are very good: i.e., probably dependent on the variational properties of the admittance parameters obtained by the Galerkin procedure.

## VII. CONCLUSION

We have presented an efficient technique for the accurate analysis of shielded multilayered printed circuits composed of arbitrarily shaped metallic areas. This technique is based on an IE solved by using the MoM with entire domain basis functions. The basis functions are efficiently evaluated by the BI-RME method. This leads to MoM matrices of small size, even in the case of complex circuits. Moreover, the transformation of the coupling integrals from surface to line integrals permits to use the basic outputs of the BI-RME method for calculating the MoM matrix.

The analyzes of circuits of practical interest have been reported and compared with both theoretical and experimental data, showing that the approach is indeed feasible and leads to a software code, which is both efficient and accurate.

## APPENDIX

Following a procedure similar to the one presented in [21], [22], the transformation of the coupling integrals from surface to line integrals is based on the application of Green's identity (see, for instance, [23, Appendix 2, eq. 45])

$$\int_{S_p} \vec{\mathcal{A}} \cdot \nabla_T \mathcal{B} dS = \int_{\partial S_p} \mathcal{B} \vec{n} \cdot \vec{\mathcal{A}} d\ell + \int_{\overline{\partial S_p}} \mathcal{B} \vec{n} \cdot \vec{\mathcal{A}} d\ell - \int_{S_p} \mathcal{B} \nabla_T \cdot \vec{\mathcal{A}} dS. \quad (\text{A.1})$$

In the case of internal patches, obviously, the port segment  $\overline{\partial S_p}$  is not defined and the line integral on  $\overline{\partial S_p}$  vanishes.

## Derivation of (20)

The electric modal fields are related to scalar potentials through (8) and (3), thus resulting in

$$\begin{aligned} \int_{S_p} \vec{e}_r^{\prime(p)} \cdot \vec{\mathcal{E}}_m dS &= \int_{S_p} \vec{u}_z \times \frac{\nabla_T \psi_r^{\prime(p)}}{\kappa_r^{\prime(p)}} \cdot \frac{\nabla_T \chi_m'}{k_m'} dS \\ &= -\frac{1}{\kappa_r^{\prime(p)} k_m'} \int_{S_p} \nabla_T \psi_r^{\prime(p)} \cdot \vec{u}_z \times \nabla_T \chi_m' dS. \end{aligned} \quad (\text{A.2})$$

By applying (A.1) to (A.2) with  $\vec{\mathcal{A}} = \vec{u}_z \times \nabla_T \chi_m'$  and  $\mathcal{B} = \psi_r^{\prime(p)}$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{\prime(p)} \cdot \vec{\mathcal{E}}_m dS &= -\frac{1}{\kappa_r^{\prime(p)} k_m'} \int_{\partial S_p} \psi_q^{\prime(p)} \vec{n}_p \cdot \vec{u}_z \times \nabla_T \chi_m' d\ell \\ &\quad - \frac{1}{\kappa_r^{\prime(p)} k_m'} \int_{\overline{\partial S_p}} \psi_q^{\prime(p)} \vec{n}_p \cdot \vec{u}_z \times \nabla_T \chi_m' d\ell \\ &\quad + \frac{1}{\kappa_r^{\prime(p)} k_m'} \int_{S_p} \psi_r^{\prime(p)} \nabla_T \cdot (\vec{u}_z \times \nabla_T \chi_m') dS. \end{aligned} \quad (\text{A.3})$$

On the right-hand side of (A.3), the surface integral vanishes because of

$$\nabla_T \cdot (\vec{u}_z \times \nabla_T \mathcal{F}) = 0 \quad (\text{A.4})$$

for any scalar function  $\mathcal{F}$  (see, for instance, [23, Appendix 2, eq. 38]), whereas the line integral on  $\partial S_p$  vanishes since  $\psi_r^{\prime(p)} = 0$  on  $\partial S_p$ . Moreover, the line integral on  $\overline{\partial S_p}$  is not defined for internal patches and, in the case of patches attached to ports, it vanishes since  $\vec{n}_p \cdot \vec{u}_z \times \nabla_T \chi_m' = -\partial \chi_m' / \partial t_p = 0$  on  $\overline{\partial S_p}$ . This proves (20).

## Derivation of (21)

From (9) and (3), we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{\prime\prime(p)} \cdot \vec{\mathcal{E}}_m dS &= \int_{S_p} \frac{\nabla_T \psi_r^{\prime\prime(p)}}{\kappa_r^{\prime\prime(p)}} \cdot \frac{\nabla_T \chi_m'}{k_m'} dS \\ &= \frac{1}{\kappa_r^{\prime\prime(p)} k_m'} \int_{S_p} \nabla_T \psi_r^{\prime\prime(p)} \cdot \nabla_T \chi_m' dS. \end{aligned} \quad (\text{A.5})$$

By applying (A.1) to (A.5) with  $\vec{\mathcal{A}} = \nabla_T \chi_m'$  and  $\mathcal{B} = \psi_r^{\prime\prime(p)}$ , and using the Helmholtz equation  $\nabla_T^2 \chi_m' = -k_m'^2 \chi_m'$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{\prime\prime(p)} \cdot \vec{\mathcal{E}}_m dS &= \frac{1}{\kappa_r^{\prime\prime(p)} k_m'} \int_{\partial S_p} \psi_r^{\prime\prime(p)} \vec{n}_p \cdot \nabla_T \chi_m' d\ell \\ &\quad + \frac{1}{\kappa_r^{\prime\prime(p)} k_m'} \int_{\overline{\partial S_p}} \psi_r^{\prime\prime(p)} \vec{n}_p \cdot \nabla_T \chi_m' d\ell \\ &\quad + \frac{k_m'}{\kappa_r^{\prime\prime(p)}} \int_{S_p} \psi_r^{\prime\prime(p)} \chi_m' dS. \end{aligned} \quad (\text{A.6})$$

By remembering that  $\psi_r^{\prime\prime(p)} = 0$  on  $\overline{\partial S_p}$ , the line integral on  $\overline{\partial S_p}$  in (A.6) vanishes.

Moreover, by applying (A.1) to (A.5) with  $\vec{\mathcal{A}} = \nabla_T \psi_r^{\prime\prime(p)}$  and  $\mathcal{B} = \chi_m'$ , and taking into account that  $\nabla_T^2 \psi_r^{\prime\prime(p)} = -\kappa_r^{\prime\prime(p)2} \psi_r^{\prime\prime(p)}$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{\prime\prime(p)} \cdot \vec{\mathcal{E}}_m dS &= \frac{1}{\kappa_r^{\prime\prime(p)} k_m'} \int_{\partial S_p} \chi_m' \vec{n}_p \cdot \nabla_T \psi_r^{\prime\prime(p)} d\ell \\ &\quad + \frac{1}{\kappa_r^{\prime\prime(p)} k_m'} \int_{\overline{\partial S_p}} \chi_m' \vec{n}_p \cdot \nabla_T \psi_r^{\prime\prime(p)} d\ell \\ &\quad + \frac{\kappa_r^{\prime\prime(p)}}{k_m'} \int_{S_p} \psi_r^{\prime\prime(p)} \chi_m' dS. \end{aligned} \quad (\text{A.7})$$

By remembering that  $\partial \psi_r^{\prime\prime(p)} / \partial n_p = 0$  on  $\partial S_p$  and that  $\chi_m' = 0$  on  $\overline{\partial S_p}$ , the line integrals in (A.7) vanish. Therefore, by substituting the surface integral on the right-hand side of (A.6) into (A.7), and considering  $\kappa_r^{\prime\prime(p)2} \neq k_m'^2$ , we finally obtain (21).

## Derivation of (22)

From (12) and (3), we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{0(p)} \cdot \vec{\mathcal{E}}_m dS &= \int_{S_p} \vec{u}_z \times \nabla_T \psi_r^{0(p)} \cdot \frac{\nabla_T \chi_m'}{k_m'} dS \\ &= \frac{1}{k_m'} \int_{S_p} \nabla_T \psi_r^{0(p)} \cdot \nabla_T \chi_m' \times \vec{u}_z dS. \end{aligned} \quad (\text{A.8})$$

By applying (A.1) to (A.8) with  $\vec{\mathcal{A}} = \nabla_T \chi_m' \times \vec{u}_z$  and  $\mathcal{B} = \psi_r^{0(p)}$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}_r^{0(p)} \cdot \vec{\mathcal{E}}_m dS &= \frac{1}{k_m'} \int_{\partial S_p} \psi_r^{0(p)} \vec{n}_p \cdot \nabla_T \chi_m' \times \vec{u}_z d\ell \\ &\quad + \frac{1}{k_m'} \int_{\overline{\partial S_p}} \psi_r^{0(p)} \vec{n}_p \cdot \nabla_T \chi_m' \times \vec{u}_z d\ell \\ &\quad - \frac{1}{k_m'} \int_{S_p} \psi_r^{0(p)} \nabla_T \cdot (\nabla_T \chi_m' \times \vec{u}_z) dS. \end{aligned} \quad (\text{A.9})$$

On the right-hand side of (A.9), the surface integral vanishes due to (A.4). With regard to the line integral on  $\partial S_p$ ,  $\psi_r^{0(p)} = \text{const}$  on each contour  $\partial S_p^i$ . In the case of a closed contour  $\partial S_p^i$

$$\int_{\partial S_p^i} \vec{n}_p \cdot \nabla_T \chi_m' \times \vec{u}_z d\ell = \int_{\partial S_p^i} \nabla_T \chi_m' \cdot \vec{t}_p d\ell = 0 \quad (\text{A.10})$$

(see [23, Appendix 2, eq. 55]). In the case of an open line  $\partial S_p^i$  (only for patches attached to ports)

$$\begin{aligned} \int_{\partial S_p^i} \vec{n}_p \cdot \nabla_T \chi'_m \times \vec{u}_z d\ell &= \int_{\partial S_p^i} \nabla_T \chi'_m \cdot \vec{t}_p d\ell \\ &= \chi'_m(P) - \chi'_m(Q) \end{aligned} \quad (\text{A.11})$$

where  $P$  and  $Q$  are extreme points of the line  $\partial S_p^i$ , which are located on the box wall  $\partial S_{\text{box}}$ , where  $\chi'_m = 0$ . Finally, the line integral on  $\overline{\partial S_p}$  vanishes since  $\vec{n}_p \cdot \nabla_T \chi'_m \times \vec{u}_z = \partial \chi'_m / \partial t_p = 0$  on  $\overline{\partial S_p}$ . This proves (22).

#### Derivation of (23)

From (8) and (4), we have

$$\begin{aligned} \int_{S_p} \vec{e}'^{(p)} \cdot \vec{\mathcal{E}}''_m dS &= \int_{S_p} \left( \vec{u}_z \times \frac{\nabla_T \psi_r^{(p)}}{\kappa_r^{(p)}} \right) \cdot \left( \frac{\nabla_T \chi''_m}{k_m''} \times \vec{u}_z \right) dS \\ &= -\frac{1}{\kappa_r^{(p)} k_m''} \int_{S_p} \nabla_T \psi_r^{(p)} \cdot \nabla_T \chi''_m dS. \end{aligned} \quad (\text{A.12})$$

Therefore, the derivation of (23) is similar to the one of (21), only taking into account the different boundary condition of the scalar potential  $\psi_r^{(p)}$ .

#### Derivation of (24)

From (9) and (4), we have

$$\begin{aligned} \int_{S_p} \vec{e}''^{(p)} \cdot \vec{\mathcal{E}}''_m dS &= \int_{S_p} \frac{\nabla_T \psi_r''^{(p)}}{\kappa_r''^{(p)}} \cdot \left( \frac{\nabla_T \chi''_m}{k_m''} \times \vec{u}_z \right) dS \\ &= -\frac{1}{\kappa_r''^{(p)} k_m''} \int_{S_p} \nabla_T \psi_r''^{(p)} \cdot \nabla_T \chi''_m \times \vec{u}_z dS. \end{aligned} \quad (\text{A.13})$$

By applying (A.1) to (A.13) with  $\vec{\mathcal{A}} = \nabla_T \chi''_m \times \vec{u}_z$  and  $\mathcal{B} = \psi_r''^{(p)}$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}''^{(p)} \cdot \vec{\mathcal{E}}''_m dS &= \frac{1}{\kappa_r''^{(p)} k_m''} \int_{\partial S_p} \psi_r''^{(p)} \vec{n}_p \cdot \nabla_T \chi''_m \times \vec{u}_z d\ell \\ &\quad + \frac{1}{\kappa_r''^{(p)} k_m''} \int_{\overline{\partial S_p}} \psi_r''^{(p)} \vec{n}_p \cdot \nabla_T \chi''_m \times \vec{u}_z d\ell \\ &\quad - \frac{1}{\kappa_r''^{(p)} k_m''} \int_{S_p} \psi_r''^{(p)} \nabla_T \cdot (\nabla_T \chi''_m \times \vec{u}_z) dS. \end{aligned} \quad (\text{A.14})$$

On the right-hand side of (A.14), the surface integral vanishes due to (A.4). In the case of patches connected to ports, the line integral on  $\overline{\partial S_p}$  vanishes since  $\psi_r''^{(p)} = 0$  on  $\overline{\partial S_p}$ . Moreover, in the line integral  $\partial S_p$ ,  $\vec{n}_p \cdot \nabla_T \chi''_m \times \vec{u}_z = \nabla_T \chi''_m \cdot \vec{t}_p$ . This proves (24).

#### Derivation of (25)

From (12) and (4), we have

$$\begin{aligned} \int_{S_p} \vec{e}^{0(p)} \cdot \vec{\mathcal{E}}''_m dS &= \int_{S_p} \left( \vec{u}_z \times \nabla_T \psi_r^{0(p)} \right) \cdot \left( \frac{\nabla_T \chi''_m}{k_m''} \times \vec{u}_z \right) dS \\ &= -\frac{1}{k_m''} \int_{S_p} \nabla_T \psi_r^{0(p)} \cdot \nabla_T \chi''_m dS. \end{aligned} \quad (\text{A.15})$$

By applying (A.1) to (A.15) with  $\vec{\mathcal{A}} = \nabla_T \psi_r^{0(p)}$  and  $\mathcal{B} = \chi''_m$ , we have

$$\begin{aligned} \int_{S_p} \vec{e}^{0(p)} \cdot \vec{\mathcal{E}}''_m dS &= \frac{1}{k_m''} \int_{\partial S_p} \chi''_m \vec{n}_p \cdot \nabla_T \psi_r^{0(p)} d\ell \\ &\quad + \frac{1}{k_m''} \int_{\overline{\partial S_p}} \chi''_m \vec{n}_p \cdot \nabla_T \psi_r^{0(p)} d\ell \\ &\quad - \frac{1}{k_m''} \int_{S_p} \chi''_m \nabla_T^2 \psi_r^{0(p)} dS. \end{aligned} \quad (\text{A.16})$$

On the right-hand side of (A.16), the surface integral vanishes because  $\nabla_T^2 \psi_r^{0(p)} = 0$ , whereas the line integral on  $\overline{\partial S_p}$  vanishes since  $\partial \psi_r^{0(p)} / \partial n_p = 0$  on  $\overline{\partial S_p}$ . This proves (25).

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