

Efficient Electromagnetic Modeling of Microstrip Structures in Multilayer Media

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Abstract—This paper presents an efficient method-of-moments solution of the mixed-potential integral equation for a general microstrip structure in multilayer media. In this method, the general forms of the spectral-domain Green's functions for multilayer media are derived first. The spatial-domain Green's functions are then obtained by the discrete complex-image method, which obviates the time-consuming numerical evaluation of the Sommerfeld integral. The Rao–Wilton–Glisson basis functions are employed to provide necessary flexibility to model arbitrary shapes. To expedite the computation of frequency response over a broad band, a reduced-order model is presented using asymptotic waveform evaluation. Numerical results of multilayer circuits and antennas are presented to show the efficiency and accuracy of this method.

Index Terms—Complex-image method, Green's function, method of moments, multilayer media, reduced-order model.

I. INTRODUCTION

A VARIETY of methods have been developed for electromagnetic modeling of microstrip structures, such as antennas, circuits, and interconnects [1]. A review of them reveals that the method-of-moments (MoM) solution of the integral equation has received intense attention. In this method, the evaluation of Green's functions and the choice of basis functions are crucial to obtaining an accurate and efficient solution. Therefore, many variations of this method have been implemented and investigated. Among them, the solution based on the mixed-potential integral equation (MPIE) using the Rao–Wilton–Glisson (RWG) basis functions [2] is an attractive approach [3]–[6] because the RWG basis function offers good flexibility to model arbitrarily shaped structures, and MPIE provides a less singular kernel compared with the electrical-field integral equation (EFIE).

On the other hand, when microstrip structures become more complicated, multilayer media are employed to allow for more versatile designs [6]–[9]. This fact necessitates the modeling to be capable of dealing with microstrip structures in multilayer media. To achieve this goal, the evaluation of Green's functions for multilayer media is crucial. As is well known, the spatial-domain Green's function is the inverse Hankel transform of the spectral-domain counterpart, which is

commonly known as the Sommerfeld integral (SI). Generally, the analytical solution of the SI is not available, and the numerical integration is time consuming since the integrand is both highly oscillating and slowly decaying. Several efficient techniques have been proposed to speed up this numerical integration [10], [11]. To further accelerate the computation, the scheme that employs the precomputation and interpolation techniques is popularly used [6], [12]. Another approach to rapidly evaluate the SI is the discrete complex-image method (DCIM) [13]–[17]. The basic idea of the DCIM is to approximate the spectral kernel by a sum of complex exponentials using the Prony method or the generalized pencil-of-function (GPOF) method [18]. The SI's can then be evaluated in closed forms via the Sommerfeld identity. This method has been extensively employed to analyze microstrip structures. However, most of the work is confined to single- or double-layered structures. As claimed in [6] and [20], it is difficult to extend the DCIM to multilayer media because of the difficulty associated with the extraction of the surface-wave contribution. However, since the surface-wave contribution is insignificant for the thin multilayer configurations and becomes appreciable only in the far-field region, it is not necessary to extract the surface wave for those cases [17], thereby rendering the DCIM possible and efficient for multilayer media. In this paper, we show that the DCIM is accurate for all the examples considered without the extraction of surface-wave contribution.

Furthermore, since the above analysis is implemented in the frequency domain, we have to repeat the calculation at each frequency to obtain the frequency response over a band of interest. This can be computationally intensive for electromagnetic devices with complicated frequency responses. For this reason, several different techniques have been proposed to characterize such devices by using the reduced-order model, such as the asymptotic waveform evaluation (AWE) [22], complex frequency hopping (CFH) [23], and Padé via Lanczos (PVL) [24]. All of them were originally developed in the circuit community. The basic idea of these techniques is to approximate the frequency response or the transfer function by a low-order rational function: the Padé approximant. In AWE and CFH, the Padé approximant of the frequency response is obtained by the moment-matching process. In PVL, the Padé approximant is obtained by the Lanczos process so that the direct calculation of moments is eluded. Recently, these techniques have been extended for electromagnetic analysis [25]–[28]. A parallel effort ongoing in the electromagnetic community is the development of the model-based parameter estimation (MBPE) [29], [30]. Instead of matching moments in

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by the reflection and transmission coefficients. One approach to simplify the derivations is to employ a transmission-line network analog of the multilayer medium, in which each layer is represented by a transmission-line section and the TE and TM waves are represented by two independent transmission lines [20], [21]. In this paper, the approach described in [19] is employed. For planar microstrip configurations, we consider a horizontal electric dipole (HED) embedded in layer m . The field point is located in layer n . The spectral-domain Green's functions for scalar and vector potentials can be written as

$$\tilde{G}_a = \frac{\mu_m}{2jk_{mz}} F_n^{\text{TE}} \quad (5)$$

$$\tilde{G}_q = \frac{1}{2j\epsilon_n k_{mz} k_\rho^2} \left(\frac{\mu_m k_n^2}{\mu_n} F_n^{\text{TE}} - jk_{mz} \frac{\partial}{\partial z} F_n^{\text{TM}} \right) \quad (6)$$

where F_n^{TE} and F_n^{TM} are the TE and TM components in layer n , respectively, which can be derived as in [19].

C. DCIM

Once the spectral-domain Green's functions are obtained, the next step is to evaluate the SI, which is very time consuming because of the highly oscillating and slowly decaying behavior of the integrand. Although some techniques have been proposed for fast evaluation of the SI, the numerical integration is still the bottleneck for the design of a fast algorithm. The DCIM obviates numerical integration and represents the SI in a closed form. Initially, the DCIM first extracts the quasi-static and surface-wave contributions from the spectral-domain kernel, then approximates the remaining kernel by a sum of complex exponentials, and finally evaluates the integral via the Sommerfeld identity. Since the extraction of the surface-wave contributions for multilayer media is cumbersome, it is difficult to extend the DCIM to multilayer media. Recently, the GPOF method has been applied to cast the Green's functions into closed forms, which is more robust and less noise sensitive compared to the original Prony method. The use of GPOF also shows that the extraction of the surface-wave contributions is not necessary for structures with insignificant surface waves. This makes it easy to employ the DCIM for multilayer media.

To apply the DCIM, we rewrite the spectral-domain Green's function in a simple form as

$$\tilde{G} = \frac{F}{2jk_{mz}}. \quad (7)$$

As the first step of the DCIM, the quasi-static contributions, which dominate as $k_\rho \rightarrow \infty$, are extracted, which makes the remaining kernel decay to zero for sufficiently large k_ρ . For the TE case, the reflection coefficients approach $(\mu_i - \mu_{i+1})/(\mu_i + \mu_{i+1})$ while, in the TM case, they tend to $(\epsilon_i - \epsilon_{i+1})/(\epsilon_i + \epsilon_{i+1})$. Therefore, we can rewrite (7) as

$$\tilde{G} = \frac{F - F_0}{2jk_{mz}} + \frac{F_0}{2jk_{mz}} \quad (8)$$

where F_0 denotes the quasi-static term. With the aid of the Sommerfeld identity, the quasi-static contribution from the

second term in (8) can be evaluated as

$$G_0 = F_0 \frac{e^{-jk_m \rho}}{4\pi\rho}. \quad (9)$$

The first term in (8) can be approximated as a sum of complex exponentials using the GPOF method. The deformed integration path in the complex k_{mz} is chosen as $k_{mz} = k_m[(1-t/T_0) - jt]$, where T_0 is determined such that $F - F_0$ becomes negligible beyond T_0 . Uniformly sampling $F - F_0$ on $t \in [0, T_0]$ and applying the GPOF method, we can approximate $F - F_0$ as

$$F - F_0 = \sum_{i=1}^{N_c} a_i e^{-k_{mz} b_i}. \quad (10)$$

Again using the Sommerfeld identity, we obtain

$$G = F_0 \frac{e^{-jk_m \rho}}{4\pi\rho} + \sum_{i=1}^{N_c} a_i \frac{e^{-jk_m r_i}}{4\pi r_i} \quad (11)$$

where $r_i = \sqrt{\rho^2 - b_i^2}$ is complex, which is the location of the complex image, and N_c is the number of complex images.

As mentioned in [15], when the source is in the bounded region of multilayer media, a modification of the DCIM is necessary. In that case, the exponentials are written in terms of the propagation coefficient of the unbounded layer instead of that of the source layer.

D. AWE Implementation

The solution can be achieved by solving (1). Given a frequency band, we have to repeat the solution of (1) at a set of frequencies to obtain the frequency response. For structures with a highly oscillatory frequency response, this process is computationally expensive. Several techniques to realize the fast frequency sweep have been developed, such as MBPE, AWE, CFH, and PVL. The basic principle of these techniques is to extract the dominant poles and residues of the frequency response and represent it by a reduced-order model, which is accomplished by constructing the Padé approximant of the frequency response.

In AWE, the solution of (1) can be expanded as a Taylor series at a given frequency point ω_0

$$I(\omega_0 + \sigma) = \sum_{n=0}^{\infty} \mathbf{M}_n \sigma^n \quad (12)$$

where \mathbf{M}_n is the n th moment vector and is given by

$$\mathbf{M}_n = \frac{1}{n!} \frac{\partial^n}{\partial \omega^n} [\bar{\mathbf{Z}}^{-1}(\omega) \mathbf{V}] \Big|_{\omega=\omega_0}. \quad (13)$$

The moments can be evaluated by the recursive relation

$$\bar{\mathbf{Z}}(\omega_0) \mathbf{M}_n = - \sum_{r=1}^n \frac{1}{r!} \frac{\partial^r}{\partial \omega^r} \bar{\mathbf{Z}}(\omega) \Big|_{\omega=\omega_0} \mathbf{M}_{n-r} \quad (14)$$

with the initial vector

$$\mathbf{M}_0 = \bar{\mathbf{Z}}^{-1}(\omega_0) \mathbf{V}. \quad (15)$$

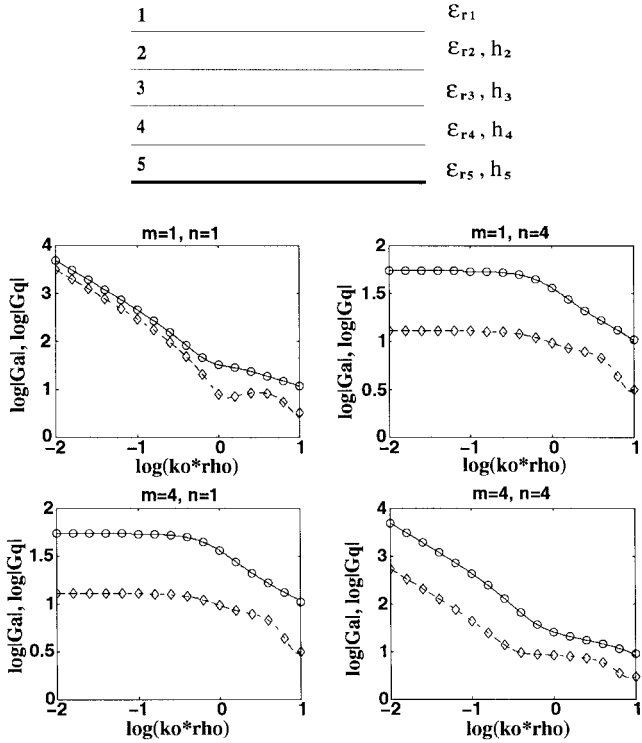


Fig. 2. Green's functions of the scalar and vector potentials in five-layer medium. The relative permittivity and thickness of each layer are as follows: $\epsilon_{r1} = 1.0$, $\epsilon_{r2} = 2.1$, $h_2 = 0.7$ mm, $\epsilon_{r3} = 12.5$, $h_3 = 0.3$ mm, $\epsilon_{r4} = 9.8$, $h_4 = 0.5$ mm, $\epsilon_{r5} = 8.6$, $h_5 = 0.3$ mm. The solid and dashed lines represent G_a and G_q , respectively, which are calculated by DCIM. The circle and diamond represent G_a and G_q , respectively, which are obtained by direct numerical integration. The frequency is 30 GHz.

To overcome the problem associated with the radius of convergence of the Taylor series expansion, the Padé approximant of $I(\omega_0 + \sigma)$ is used, which can be written as

$$I(\omega_0 + \sigma) = \frac{\mathbf{a}_0 + \mathbf{a}_1\sigma + \dots + \mathbf{a}_p\sigma^p}{1 + \mathbf{b}_1\sigma + \dots + \mathbf{b}_q\sigma^q} \quad (16)$$

where typically $p = q-1$ and the $p+q+1$ unknown coefficients of the numerator and denominator are uniquely determined by matching the first $p+q+1$ moments of (12), as is done in [25] and [26]. Once we obtain the Padé approximant, the frequency response over a frequency band can be rapidly evaluated.

Therefore, the first step of AWE is to evaluate the derivative matrices at the expansion point. As can be seen from (2), the impedance matrix is the function of frequency through the Green's functions. Unlike in [30], the Green's functions in this method have the simple expression e^{-jkr}/r , as is evident in (11), due to the use of the DCIM. Therefore, here, the derivative matrices can be simply formulated. The moment vectors can then be obtained recursively by (14). After the moment-matching process, the Padé approximant for each element can be obtained. Therefore, the frequency response over a band near the expansion point is obtained.

In many practical problems with a broad-band response, one expansion point is not sufficient to cover the entire bandwidth. In such cases, multiple expansion points are necessary. Here, a simple binary search algorithm, as described in [25] and [26], is employed to automatically choose the expansion points. As-

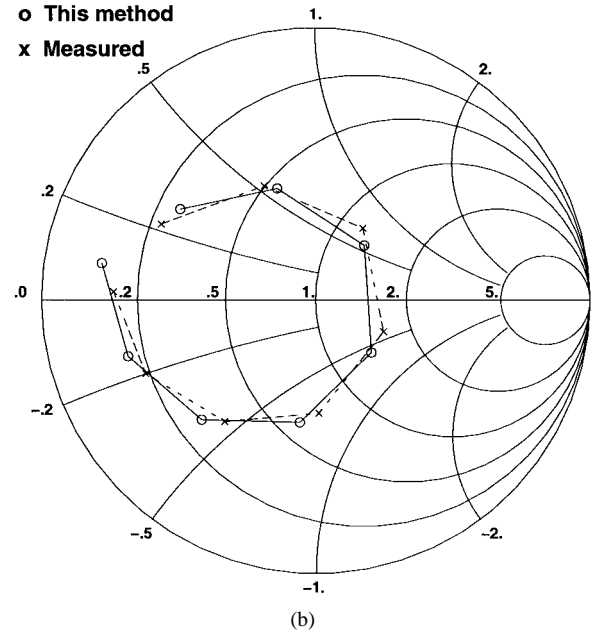
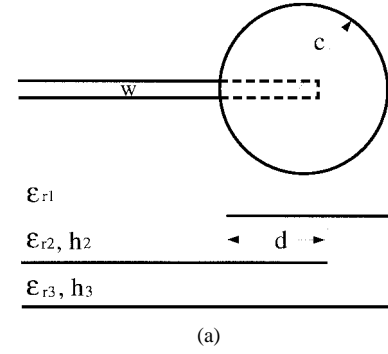


Fig. 3. Input impedance for a proximity-coupled microstrip antenna. $\epsilon_{r1} = 1.0$, $\epsilon_{r2} = \epsilon_{r3} = 2.62$, $h_2 = h_3 = 1.59$ mm, $w = 4.372$ mm, $c = 17.5$ mm, $d = 17.5$ mm. The frequency is from 2.80 to 3.15 GHz with an increment of 0.05 GHz.

sume that we are interested in the frequency response over the band $[f_{\min}, f_{\max}]$. We first compute the reduced-order models using f_{\min} and f_{\max} as the expansion points. We then compute the frequency response over the entire band $[f_{\min}, f_{\max}]$. If the two responses are within an acceptable error tolerance, the procedure has converged. If not, an additional reduced-order model is then computed at a new expansion point $(f_{\min} + f_{\max})/2$, and the error checking process is repeated in the two new subintervals. This process is repeated until the two reduced-order models bordering all subintervals give the frequency responses within the prescribed error tolerance.

It is known that AWE has the problem of instability in the computation of the Padé approximation due to the ill-conditioned moment-matching process [24]. However, this is not a problem in our method because the order of the Padé approximant is typically chosen as $q \leq 8$, for which the instability has never been observed.

III. NUMERICAL RESULTS

We first verify the spatial-domain Green's functions obtained by the DCIM by considering a five-layer medium. The

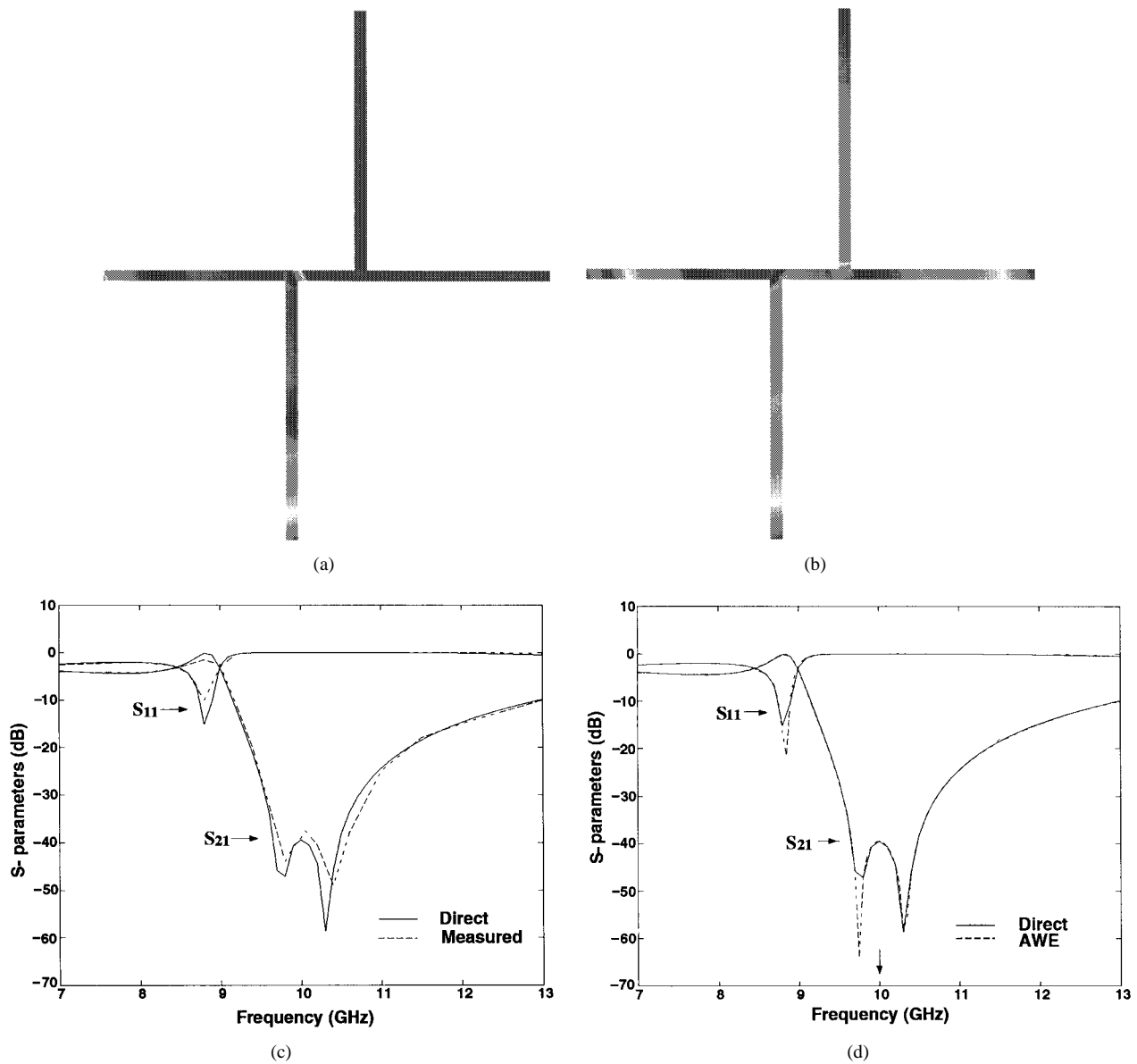


Fig. 4. Current distributions and S -parameters for a microstrip double stub. $\epsilon_r = 9.9$, $h = 0.127$ mm. The linewidth is 0.122 mm. The stub length is 2.921 mm. The spacing between two stubs is 0.757 mm. (a) $f = 10$ GHz. (b) $f = 13$ GHz.

TABLE I
COMPARISON OF THE CPU TIME USING THE DIRECT CALCULATION AND AWE

	Direct			AWE			Speed-up
	T_{freq} (sec)	N_{freq}	T_1 (sec)	T_{expn} (sec)	N_{expn}	T_2 (sec)	
Ex. 1	6.4	60	384.0	18.1	1	18.1	21.2
Ex. 2	104.0	80	8320.0	350.2	3	1050.6	7.9
Ex. 3	27.0	400	10800.0	95.9	9	863.1	12.5

- T_{freq} : CPU time per frequency in the direct calculation;
- N_{freq} : Number of frequencies in the direct calculation;
- T_1 : Total CPU time by the direct calculation;
- T_{expn} : CPU time per expansion point with AWE;
- N_{expn} : Number of expansion points with AWE;
- T_2 : Total CPU time with AWE.

first layer is free space. The fifth layer is backed by a ground plane. The frequency is 30 GHz. We seek Green's functions for the four cases. The HED can be at the interface of the first and second layer $m = 1$ or the fourth and fifth layer $m = 4$.

The field point also has two choices of $n = 1$ or $n = 4$. The Green's functions of all these four cases are given in Fig. 2 compared to those obtained by the numerical integration along the Sommerfeld integration path (SIP). A good agreement can be observed. In this calculation, we choose $T_0 = 20$ and $N_c = 5$. The saving of the computation time is very significant. For instance, the average central processing unit (CPU) time to obtain one Green's function is 15 s using the numerical integration approach. In contrast, the DCIM only takes less than 0.2 s. The computation is performed on a 266-MHz DEC Alpha workstation. Furthermore, the numerical integration is performed from point to point; however, the DCIM is a plane-to-plane calculation so that the source and field points on the same z -plane share the same DCI's, which further enhances the efficiency of the DCIM.

After the validation of this technique, we incorporate the DCIM into the MoM solution of MPIE for microstrip structures in multilayer media. The example we consider is a

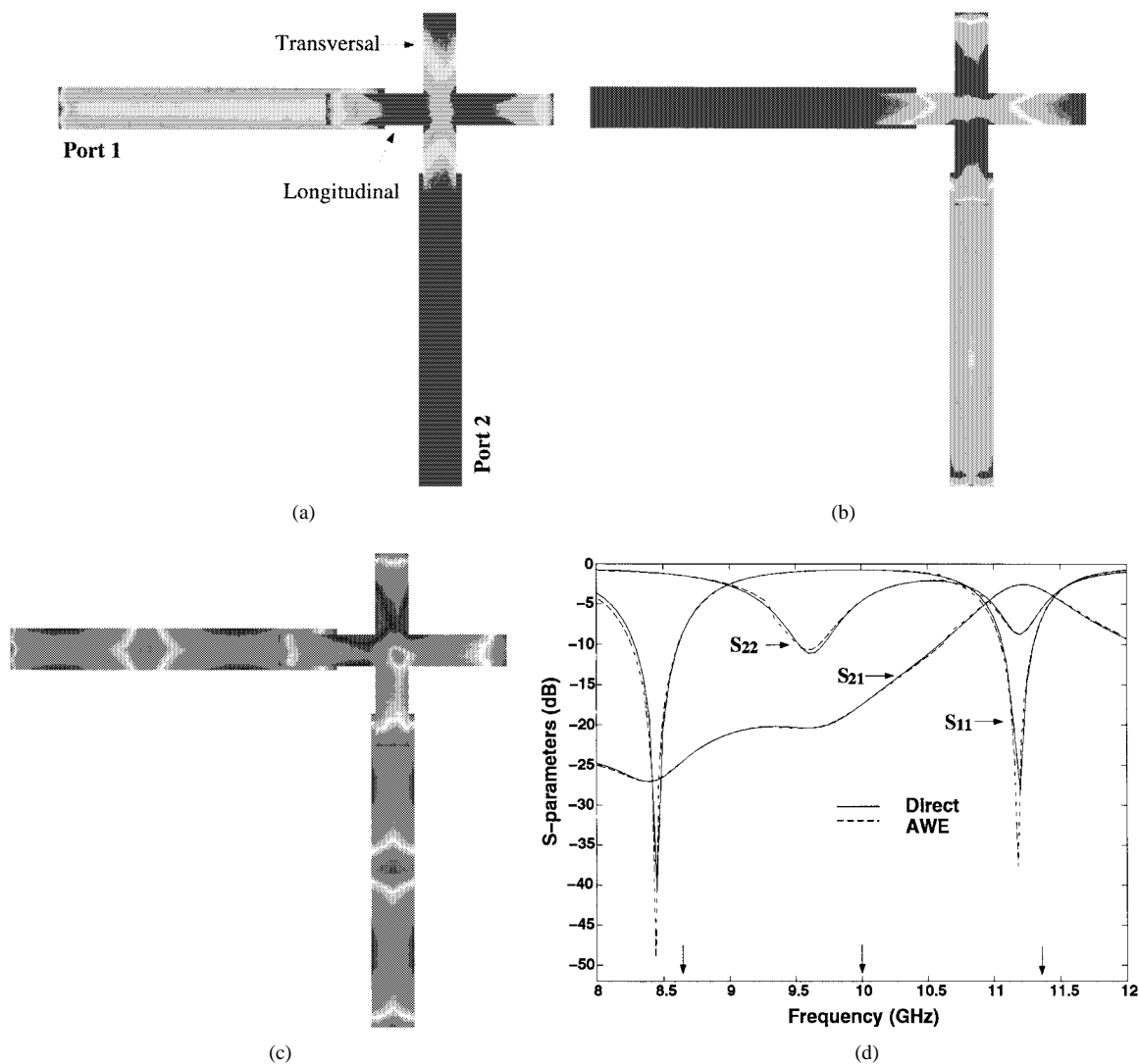


Fig. 5. Current distributions and S -parameters for a two-port asymmetric antenna. (a) $f = 8.4$ GHz. (b) $f = 9.6$ GHz. (c) $f = 11.3$ GHz.

proximity-coupled circular patch antenna [31]. The two substrate layers have the same relative permittivity $\epsilon_r = 2.62$ and thickness 1.59 mm. The microstrip line is on the bottom layer and the patch is on the top layer. The radius of the patch is 17.5 mm and the width of the microstrip line is 4.372 mm. The overlapping length may be used to control the coupling. Here, it is 17.5 mm. The number of unknowns is 686. The CPU time for one frequency is 89.3 s. The input impedance with the reference plane 79.0 mm away from the patch center is given in Fig. 3, which agrees well with the measured data from [31].

In the following three examples, AWE is applied to expedite the calculation of frequency response over a bandwidth. In all the cases, we use $q = 8$. The CPU time using AWE is compared with that employing the direct calculation to demonstrate the efficiency, as shown in Table I. We first consider a microstrip double-stub on a single layer, which has the relative permittivity $\epsilon_r = 9.9$ and the thickness 0.127 mm [12], [32]. The number of unknowns is 205. Current distributions at two different frequencies are shown in Fig. 4(a) and (b). The direct calculation gives a very good result

comparing with the measured data [12], [32], as shown in Fig. 4(c). In this calculation, the CPU time at each frequency is 6.4 s and totally 60 frequencies are sampled to obtain the accurate results. Therefore, the total CPU time is 384 s. Using AWE and choosing only one expansion point $f = 10.0$ GHz, we can obtain a very good result, as given by Fig. 4(d). In the calculation with AWE, the CPU time is 18.1 s. Thus, using AWE is approximately 21 times faster. Note that the sampling points are not dense enough to catch the null at approximately 9.75 GHz for the direct calculation.

We now consider a two-port asymmetric antenna [33]. There are two orthogonally crossed dipoles on the top layer. They have different lengths so that dual-frequency operation is achieved. The transversal one has a length of 11.9 mm and the longitudinal one has a length of 10.2 mm. The width of both is 1.7 mm. The two substrate layers have the same relative permittivity $\epsilon_r = 2.17$. The top layer has a thickness of 1.6 mm and the bottom layer has a thickness of 0.8 mm. On the bottom layer, the feeding lines are 2.2-mm wide. Fed from port 1, the longitudinal dipole is excited at the resonant frequency of 8.4 GHz. The current distribution is given in

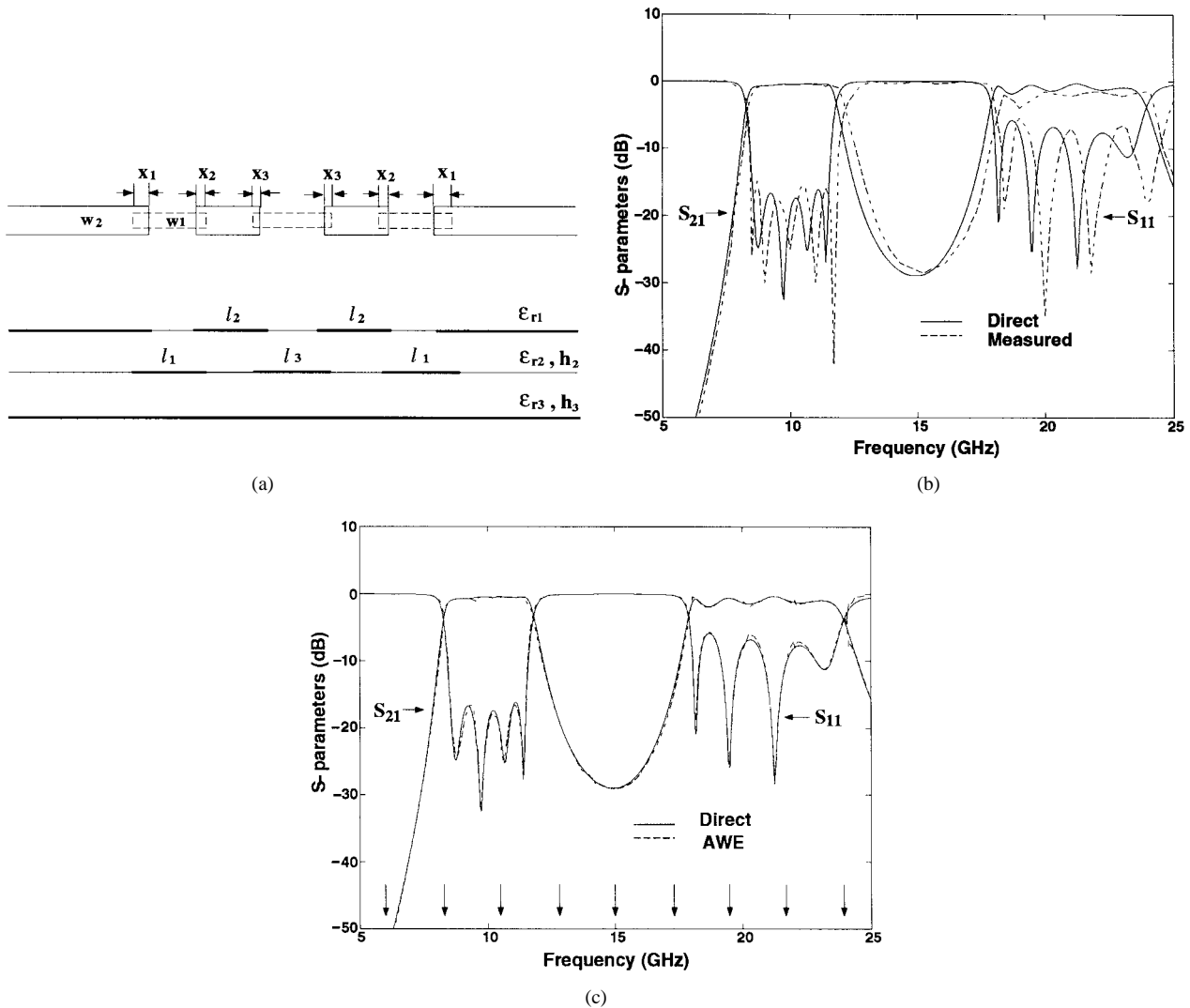


Fig. 6. S -parameters for an overlap-gap-coupled microstrip filter. $\epsilon_{r1} = 1.0$, $\epsilon_{r2} = 9.8$, $\epsilon_{r3} = 2.2$, $h_2 = h_3 = 0.01$, $l_1 = 0.2752$, $l_2 = 0.2542$, $l_3 = 0.2851$, $x_1 = 0.0516$, $x_2 = 0.0152$, $x_3 = 0.0106$. Unit of dimension: inch.

Fig. 5(a). Fed from port 2, the transversal dipole is excited at its resonant frequency of 9.6 GHz. The current is shown in Fig. 5(b). At 11.3 GHz, the coupling bend consisting of two perpendicular half-dipole is resonant. The incident power is essentially transmitted from one port to the other, and the current distribution is shown in Fig. 5(c). The magnitude of S -parameters is given in Fig. 5(d). The measured data are available from [33]. The good agreement can be observed. In this example, the number of unknowns is 912. The direct method takes 104 s for each frequency, and a total of 80 sampling frequencies are used to obtain the accurate results. Therefore, the total CPU time is approximately 139 min. In contrast, with AWE, only three expansion frequencies are needed. At each expansion point, the CPU time is 350.2 s. Totally, the CPU time is 17.5 min. Thus, using AWE is approximately eight times faster.

The final example is an overlap-gap-coupled microstrip filter [6], [34], illustrated in Fig. 6(a), which is a two-layer geometry. The top layer is alumina with a relative permittivity of 9.8 and a thickness of 0.254 mm. The bottom layer is duroid with a relative permittivity of 2.2 and a thickness

of 0.254 mm. The geometrical parameters are the same as those in [34]. The number of unknowns is 503. In the direct calculation, the CPU time at each frequency is 27 s, and a total of 400 sampling frequencies are used to obtain the accurate results. Therefore, the total CPU time is approximately 3 h. The computed S -parameters in comparison with the measured data are shown in Fig. 6(b). The small discrepancy with the measured data is due to the finite thickness of metallization and the manufacturer's tolerance of substrates, as discussed in [6] and [34]. In contrast, with AWE, only nine expansion frequencies are chosen automatically by the binary search. At each expansion point, the CPU time is 95.9 s. In total, the CPU time is 863.1 s. Thus, the method with AWE is approximately 12.5 times faster than the direct calculation. The results without and with AWE are in excellent agreement, as shown in Fig. 6(c).

IV. CONCLUSION

This paper presents an efficient MoM solution of the MPIE formulation for a general microstrip structure in a multilayer

medium. In this method, the spectral-domain Green's functions for multilayer media are derived and the spatial-domain Green's functions in the form of SI's are then evaluated by the DCIM, which obviates the time-consuming numerical integration. The RWG basis functions are employed to provide necessary flexibility to model arbitrary shapes. The AWE is applied to efficiently evaluate the frequency response over a band of interest. Some multilayer circuits and antennas are analyzed to demonstrate the efficiency and accuracy of this method. We remark that since the surface-wave contribution is not extracted, the error of the DCIM employed in this paper increases as the $k_0\rho$ increases. For all the examples, the error is within an acceptable tolerance for $k_0\rho < 100$ or $\rho < 16\lambda_0$, where λ_0 is the free-space wavelength. Most microstrip circuit components have a dimension much smaller than $16\lambda_0$, which is the primary reason for the accuracy of the proposed method.

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