

A Quasi-One-Dimensional Integration Technique for the Analysis of Planar Microstrip Circuits via MPIE/MoM

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Abstract—The mixed-potential integral-equation approach, using spatial-domain closed-form Green's functions, and discretized with the method-of-moments, is a state-of-the-art method for the analysis of planar microstrip circuits. One of its most time-demanding tasks is the evaluation of the impedance matrix terms, which typically requires the numerical computation of two-dimensional integrals. A method based on suitable changes of coordinates and domains is introduced in this paper in order to reduce such integrals to a quasi-one-dimensional numerical integration, with a substantial enhancement in the efficiency of the analysis, without affecting the accuracy of the approach. Results are given demonstrating, for practical accuracy values, an improvement of typically one order of magnitude in simulation times.

I. INTRODUCTION

EFFICIENT modeling of printed circuits and antennas is crucial in current microwave engineering [1], [2] and has stimulated several contributions. In particular, a mixed-potential integral equation (MPIE) was proposed by Mosig [3], [4]. More recently, the latter formulation was enhanced by the introduction of suitable closed-form spatial-domain Green's functions [5] and suitable transformations of the impedance matrix [6]. Considerable efforts are currently made in order to improve the efficiency and accuracy of these numerical methods, such as the inclusion of three-dimensional (3-D) unknown currents, efficient choices for the Sommerfeld integration paths [7], and the enhancement of the complex-image method [8], [9] for multi-level stratified microstrip lines [10], [11]. The above-mentioned contributions are mainly in the direction of reducing the computation time required for filling the impedance matrix. Also on this subject, and more recently, papers have been proposed that discuss a clever analysis of basis functions behavior [12], with a significant improvement of space and spectral integrations of coupling integrals [13]–[19].

The numerical core of method-of-moments (MoM) approaches for the analysis of microstrip circuits is represented by both the computation of the impedance matrix and the solution of the corresponding linear system. The computation effort for the evaluation of the impedance matrix is basically determined by the numerical evaluation of some reaction integrals. Therefore, their efficient and accurate solution,

possibly appropriate for a wide class of Green's functions, is of paramount importance.

In this paper, we propose an efficient method for evaluating the impedance matrix elements of the circuit via a suitable computation of the relevant reaction integrals. The two-dimensional (2-D) numerical integration encountered in previous approaches [20] is reduced to a quasi-one-dimensional (1-D) numerical integration. Even though this is achieved by partitioning the problem domain into equal cells, this does not represent a practical limitation to the presented technique, as different optimum sizes for the elementary cells can be used in several regions of a single circuit so that an optimum accuracy is ensured. The method proposed is valid for a very large class of Green's function (the only hypothesis is that the Green's function depends only on the source–test distance), and can also be applied to circuits with more than one dielectric layer.

This paper is organized as follows. First, for the reader's convenience, the MPIE formulation is resumed; afterwards, the quasi-1-D integration method is described, some results are given, and finally, conclusions are drawn.

II. ELECTRIC-FIELD MPIE WITH CLOSED-FORM GREEN'S FUNCTIONS

We consider N -port planar circuits with infinite transverse dimensions for both the dielectric and ground plane; the metalization thickness is assumed negligible. In order to achieve improved convergence properties, we select the MPIE formulation [3], [4], which is solved by considering closed-form Green's functions in the spatial domain and by using the MoM.

Spatial-domain mixed —potential Green's functions for a layered medium are expressed by Sommerfeld integrals [21] whose integrands are slowly decaying oscillating functions, hence, the calculation is very time consuming. A possible approach to circumvent this problem is the quasi-dynamic image model [22], which is not accurate enough when surface and leaky wave effects must be accounted for [23].

The evaluation of the above-mentioned Green's functions in closed form is performed as suggested in [5], [9], and [24]. Accordingly, the spatial-domain mixed-potential Green's functions are written in the following manner:

$$\begin{aligned} G_{xx}^A &= G_{xx,0}^A + G_{xx,SW}^A + G_{xx,ci}^A \\ G^q &= G_0^q + G_{SW}^q + G_{ci}^q \end{aligned} \quad (1)$$

i.e., as the sum of direct terms and quasi-dynamic images ($G_{xx,0}^A$, G_0^q), surface waves ($G_{xx,SW}^A$, G_{SW}^q), and complex

Manuscript received August 8, 2000.

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Publisher Item Identifier S 0018-9480(01)01687-8.

images ($G_{xx,ci}^A$, G_{ci}^A). As well known, the complex image method is not so accurate when the source–test distance is larger than a certain threshold. This is taken into account by performing a phenomenological analysis, as described in [25]. The same analysis is also useful to cope with some singularity problems, often encountered in space-domain formulations [26].

The Galerkin's MoM is used to discretize the relevant equations, by selecting rooftop functions defined over elementary rectangular domains. This way, a linear system of size N is derived from the MPIE

$$\begin{bmatrix} Z_{xx} & Z_{xy} \\ Z_{yx} & Z_{yy} \end{bmatrix} \begin{bmatrix} I_x \\ I_y \end{bmatrix} = \begin{bmatrix} V_x \\ V_y \end{bmatrix}. \quad (2)$$

The entry Z_{ij} in the impedance matrix \mathbf{Z} is expressed by a fourfold integral, in the spatial variables x' , y' —corresponding to the source coordinates—and x , y —corresponding to the test coordinates. Part of its evaluation can be performed analytically [27] and, by paying attention to the choice of appropriate basis functions, the integrals “can be reduced to double integrals over finite domains” [20]. The unknowns I_x and I_y are the (complex) amplitudes of the basis functions. The right-hand-side (RHS) vector $[V]$ depends on the excitation applied to the microstrip network.

III. QUASI-1-D INTEGRATION

In order to compute the impedance matrix, a time-demanding 2-D numerical integration was employed in earlier methods, even though several efforts have been made to improve the convergence of this computation [27], [20]. Some attempts in this sense have also been made by focusing on appropriate choices of the basis and test functions [28], with attention paid on the meshing performed on the problem's domain. In this last approach [28], the entries of the matrix in (2) are expressed as a four-dimensional (4-D) integral

$$\begin{aligned} \langle f, G * g \rangle &= \int_{D_x} \int_{D_y} f(x, y) \int_{D_{x'}} \int_{D_{y'}} G(r) g(x', y') dx' dy' dx dy \\ &\quad (3) \end{aligned}$$

where D_x , D_y , $D_{x'}$, and $D_{y'}$ are the domains along x , y , x' , and y' of the interacting cells (Fig. 1), $r = \sqrt{(x - x')^2 + (y - y')^2}$, and the fourfold integration appearing in (3) is therein transformed into a 2-D one.

In this paper, we show how integration (3) can be reduced to a quasi-1-D integral. The basic observations are: 1) the dependence of the Green's function only on the source–test distance r ; 2) the possibility of reducing the remaining terms in the integration kernel (3) to functions with the same behavior; and 3) the possibility of numerically evaluating the contour of the integration domains for the above-mentioned remaining terms. The three observations lead to the following procedure: first, the four integrals can be considered as a 2-D convolution, which can be analytically solved inside each elementary cell of the circuit. This way, (3) is reduced to a 2-D integration, whose kernel is the

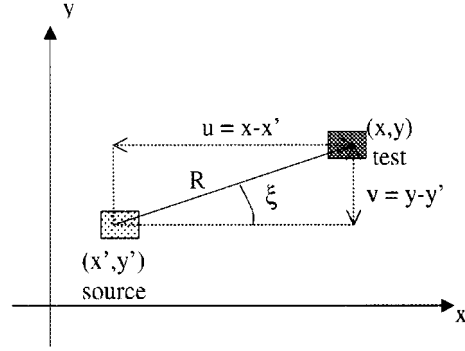


Fig. 1. Reference system and the relative changes of coordinates. One source and one test current cell are sketched.

product between the Green's function and the function attained from the 2-D convolution. A final change of coordinates is now sufficient to separate the two integrals so that one integration is numerically performed with high efficiency by evaluating an analytical function in a small number of points (less than ten). Hence, only a 1-D integration must be performed to completely evaluate (3).

We continue now to describe with more details the proposed integration technique. The usual change of variables

$$\begin{aligned} x - x' &= u \\ x + x' &= p \\ y - y' &= v \\ y + y' &= q \end{aligned} \quad (4)$$

reduces the problem to a double integration, hence, providing

$$\begin{aligned} \iint f(x, y) \iint G(\sqrt{(x - x')^2 + (y - y')^2}) &\cdot g(x', y') dx' dy' dx dy \\ &= \frac{1}{4} \iint G(u, v) \iint f\left(\frac{u+p}{2}, \frac{v+q}{2}\right) \\ &\cdot g\left(\frac{p-u}{2}, \frac{q-v}{2}\right) dp dq du dv \end{aligned} \quad (5)$$

(variables x , x' , y , y' , u , and v are introduced and described in Fig. 1).

Letting D_x be the domain for the x -variable, we can cast

$$\begin{aligned} S(u, v) &\triangleq \frac{1}{4} \int_{D(p)} \int_{D(q)} f\left(\frac{u+p}{2}, \frac{v+q}{2}\right) \\ &\cdot g\left(\frac{p-u}{2}, \frac{q-v}{2}\right) dp dq. \end{aligned} \quad (6)$$

With a suitable change of coordinates (see the Appendix for details), the integration (6) can be transformed into a bidimensional convolution

$$S(u, v) = S(\zeta, \eta) = \int_{D(\zeta)} \int_{D(\eta)} f(\zeta, \eta) g(\zeta - u, \eta - v) d\zeta d\eta \quad (7)$$

which can be solved very efficiently in analytical form, as described in the Appendix. This way, (3) is transformed into a two-variable integration

$$\int_{D(u)} \int_{D(v)} G(u, v) S(u, v) du dv. \quad (8)$$

With the following new change of variables:

$$u = r \cos \xi \quad v = r \sin \xi \quad (9)$$

we can write

$$\begin{aligned} & \int_{D(u)} \int_{D(v)} G(u, v) S(u, v) du dv \\ &= \int_{r_1}^{r_2} G(r) r \int_{\xi_1(r)}^{\xi_2(r)} S(r \cos \xi, r \sin \xi) d\xi dr. \end{aligned} \quad (10)$$

Thanks to the fact that the Green's functions only depend on the source-test distance r , a function $W(r)$

$$W(r) \triangleq \int_{\xi_1(r)}^{\xi_2(r)} S(r \cos \xi, r \sin \xi) d\xi \quad (11)$$

can be evaluated nearly completely in closed form, with a very high efficiency ($\xi_1(r)$ and $\xi_2(r)$ can be numerically evaluated for each fixed value of r). Finally, we have

$$\begin{aligned} & \iint f(x, y) \iint G(\sqrt{(x-x')^2 + (y-y')^2}) \\ & \quad \cdot g(x', y') dx' dy' dx dy \\ &= \int_{r_1}^{r_2} W(r) G(r) r dr. \end{aligned} \quad (12)$$

With a suitable choice [28] of basis and test functions, $W(r)$ is integrable in the Riemann sense. The $S(u, v)$ and $W(r)$, as well as the forms of domains $D(p)$ and $D(q)$, $\xi_1(r)$ and $\xi_2(r)$, and r_1 and r_2 depend on the choice of the basis and test functions and their domain of definition.

After evaluating $W(r)$, the Z -matrix terms can be written as

$$\begin{aligned} Z_{xx} &= \int \left[W_{1x}(r) G_{xx}^A(r) - \frac{1}{\omega^2} W_{2x}(r) G^q(r) \right] r dr \\ Z_{xy} &= \int \left[-\frac{1}{\omega^2} W_{3x}(r) G^q(r) \right] r dr \\ Z_{yx} &= \int \left[-\frac{1}{\omega^2} W_{3y}(r) G^q(r) \right] r dr \\ Z_{yy} &= \int \left[W_{1y}(r) G_{yy}^A(r) - \frac{1}{\omega^2} W_{2y}(r) G^q(r) \right] r dr \end{aligned} \quad (13)$$

thus demonstrating that the elements of the impedance matrix can be evaluated by solving a quasi-1-D integral.

The proposed quasi-1-D formulation can be exploited when equal cells are used in a rectangular mesh. An appropriate partitioning of the circuit into subregions can be generally used to analyze every subregion with an optimum cell size. In the Appendix, details are given about the derivation of the coefficients $W_{\{1,2,3\}\{x,y\}}$ for the case of rooftop functions.

As already mentioned, the quasi-1-D approach has been tested here in the case of a Galerkin MoM using rooftop basis and test functions. In the Appendix, the detailed formulation

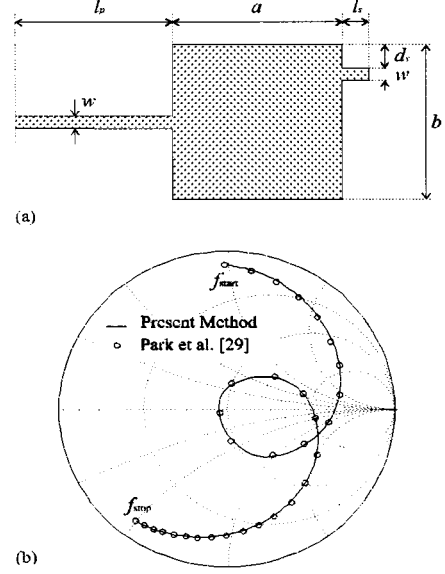


Fig. 2. (b) Input impedance on the Smith chart of: (a) the microstrip-line center-fed square antenna with a tuning stub. Physical dimensions: $\epsilon_r = 2.62$, $d = 0.794$ mm, $a = b = 28.6$ mm, $w = 2.2$ mm, $l_p = 26.4$ mm, $l_s = 4.4$ mm, $d_s = 4.4$ mm. Frequency range: 2.98–3.3 GHz. A mesh with a 3-mm edge was used for the simulations.

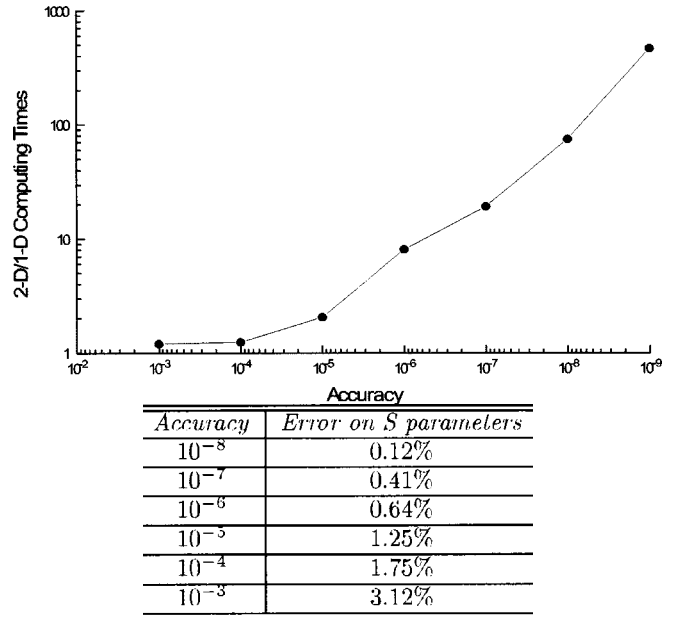


Fig. 3. Comparison between the 1-D and 2-D integration's performance. Normalized times refer to the analysis of the patch antenna in Fig. 2. They are attained as the ratio between the simulation time of the 2-D and 1-D implementation. On the x -axis, the accuracy required to make the integration converge is reported. An accuracy of 10^{-6} is typically selected for practical computation (refer to table).

for this case is given. Nevertheless, it can also be extended to other functions, such as pulse functions, attaining different forms for (13).

IV. RESULTS

The accuracy and efficiency of the implemented method is demonstrated for a patch antenna, reported in the literature and sketched in Fig. 2. In Fig. 3, we compare the time performance

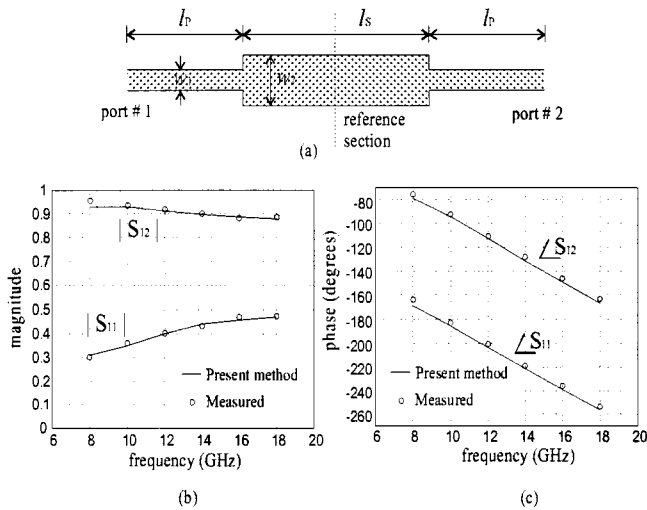


Fig. 4. (a) Scattering parameters of matching section: (b) in magnitude and (c) in phase. Physical dimensions: $\epsilon_r = 9.9$, $d = 10$ mil, $w_1 = 9.2$ mil, $w_2 = 23$ mil, $l_p = 30$ mil, $l_s = 50.6$ mil. A mesh with a 0.5-mm edge was used for the simulations.

of an MPIE/MoM package implementing a quasi-1-D integration of (3) with respect to the previous 2-D integration proposed by [28]. In the x -axis, the required numerical accuracy is reported, on the y -axis the corresponding normalized computing time for the 2-D implementation with respect to the quasi-1-D one. We define accuracy as the threshold considered to identify the integration convergence. A typical value, guaranteeing a good tradeoff between performance and accuracy is 10^{-6} (errors for scattering parameters are around 1%).

The case of Fig. 2, simulated with a mesh with a 3-mm edge, generates a matrix of dimension 240. When an accuracy of 10^{-6} is considered, the computing time for one frequency point is 2.67 s on a PC Pentium 200 MHz with the 2-D implementation, and 0.33 s using the quasi-1-D integration. A sparse banded approach is used for the system solution step, as described in [6]: it guarantees a high efficiency in the system solution time, taking advantage from the use of reordering techniques.

As easily predictable, the 1-D integration is highly superior. The numerical complexity on the number of integration points is quadratical in the case of the 2-D integration, and linear in the case of quasi-1-D solution. This is in clear accordance with the response of the reported curve.

Another evidence is reported in Fig. 4, where a two-port circuit is analyzed. The advantage of the 1-D integration is also apparent in this case (normalized 2-D/1-D simulation times are reported in Fig. 5). The case of Fig. 4, simulated with a mesh with a 0.5-mm edge, generates a matrix of dimension 416. For an accuracy of 10^{-6} , the computing time for one frequency point is 5.56 s on a PC Pentium 200 MHz with the 2-D implementation, and 0.7 s using the quasi-1-D integration. The sparse banded approach mentioned before is used [6].

The system generation (i.e., the evaluation of the entries in the impedance matrix) is a heavy computational task, as resumed in Table I. For the two circuits (the patch antenna and the two-port circuit), the percentage of time in the two main tasks (system generation and system solution) is reported. As can be seen, the system generation plays an important role. This explains the

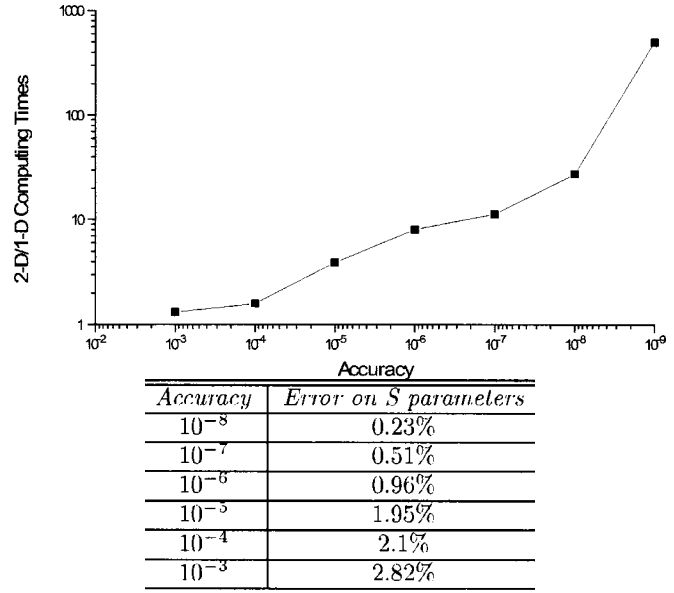


Fig. 5. As in Fig. 3, referring to the circuit in Fig. 4.

TABLE I
EFFORT REQUIRED BY THE SYSTEM GENERATION, WITH RESPECT TO THE SYSTEM SOLUTION AND THE REMAINING TASKS, FOR THE TWO ADDRESSED CASES

Circuit	System gen.	System sol.	Other
Patch Antenna	72%	16%	12%
Two-port Circuit	61%	23%	16%

importance of an efficient solution of (3), as well as the huge speed up achieved with the quasi-1-D integration.

Of course, the reported speed ups are attained when comparing the 2-D implementation with the quasi-1-D implementation of the MPIE/MoM formulation adopted in this paper. Several alternative schemes for the MPIE/MoM are available in the literature. In such cases, the applicability of the proposed strategy should be specifically investigated.

V. CONCLUSIONS

In this paper, a substantial enhancement is achieved for the efficient numerical analysis of planar microstrip circuits. A method is described based on suitable analytical changes of coordinates and domains so that the elements are evaluated by solving quasi-1-D numerical integrations instead of the 2-D integrations performed in previous approaches. The above technique, although very general, has been presented in conjunction with an MPIE formulation of the problem, and closed-form spatial-domain Green's function expressions. In such a case, the efficient evaluation of the entries of the impedance matrix is of paramount importance to achieve high performance. Results demonstrate that the proposed approach guarantees a high accuracy and decreases simulation times, in practical cases, typically of one order of magnitude.

APPENDIX

We describe the case for the xx terms of the impedance matrix; similar formulations hold for the remaining xy , yx , and yy

terms. We refer to the case of basis and test rooftop functions, i.e.,

$$J_{nx} = \begin{cases} \left(1 - \frac{|x - x_n|}{h_x}\right), & \text{if } |x - x_n| \leq h_x, |y - y_n| \leq h_y/2 \\ 0, & \text{elsewhere} \end{cases} \quad (14)$$

where $2h_x h_y$ is the surface of the current cell where the function is defined and (x_n, y_n) are the coordinates of the domain's center. Similar expressions hold for the J_{ny} . In the case of rooftop test and basis functions and xx interaction, as seen from (13), we must basically derive two functions W_{1x} and W_{2x} . In fact, the terms $Z_{xxm,n}$ can be expressed as [28]

$$Z_{xxm,n} = \langle J_{xm}, G_{xx}^A * J_{xn} \rangle - \frac{1}{\omega^2} \left\langle \frac{\partial J_{xm}}{\partial x}, G^q * \frac{\partial J_{xn}}{\partial x} \right\rangle. \quad (15)$$

It can be demonstrated that

$$\langle J_{xm}, G_{xx}^A * J_{xm} \rangle = \int_{r_1}^{r_2} W_{1x}(r) G_{xx}^A(r) r dr \quad (16)$$

$$\left\langle \frac{\partial J_{xm}}{\partial x}, G^q * \frac{\partial J_{xn}}{\partial x} \right\rangle = -\frac{1}{\omega^2} \int_{r_1}^{r_2} W_{2x}(r) G_{xx}^q(r) r dr. \quad (17)$$

Now, in accordance with (11), we indicate with S_{1x} and S_{2x} two functions so that

$$W_{1x}(r) = \int_{\xi_1}^{\xi_2} S_{1x}(r, \xi) d\xi \quad (18)$$

$$W_{2x}(r) = \int_{\xi_1}^{\xi_2} S_{2x}(r, \xi) d\xi. \quad (19)$$

Therefore, in order to find out W_{1x} and W_{2x} , we must derive the two functions S_{1x} and S_{2x} .

We first concentrate on S_{1x} . Referring to Fig. 6, (3) can be written in the following form, in the case of interaction along the x -axis for both test and basis functions, i.e., for the Z_{xxAB} terms in the impedance matrix:

$$\int_{D(Ay)} \int_{D(Ax)} J_{Ax} \int_{D(By)} \int_{D(Bx)} G(r) J_{Bx} dx' dy' dx dy \quad (20)$$

with

$$\begin{aligned} x_A - h_x &\leq x \leq x_A + h_x \\ x_B - h_x &\leq x' \leq x_B + h_x \\ y_A - h_y/2 &\leq y \leq y_A + h_y/2 \\ y_B - h_y/2 &\leq y' \leq y_B + h_y/2. \end{aligned} \quad (21)$$

By using the change of coordinates (4), we have

$$S_{1x}(u, v) \triangleq \frac{1}{4} \int_{D(p)} \int_{D(q)} J_{Apu} \left(\frac{u+p}{2}, \frac{v+q}{2} \right) \cdot \text{cdot} J_{Bpu} \left(\frac{p-u}{2}, \frac{q-v}{2} \right) dp dq. \quad (22)$$

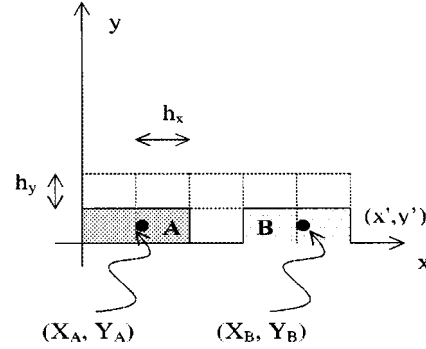


Fig. 6. Simple example for the xx interaction. Two current cells (A and B) are shown, with the relative centers.

Now, by casting

$$\zeta = \frac{\frac{p+u}{2} - x_A}{h_x} \quad \eta = \frac{\frac{p+u}{2} - x_A}{h_x} \quad (23)$$

we have $dp = 2h_x d\zeta$, $dq = 2h_y d\eta$, and

$$\frac{\frac{p-u}{2} - x_B}{h_x} = \zeta - \frac{u - (x_A - x_B)}{h_x} \quad (24)$$

$$\frac{\frac{q-v}{2} - y_B}{h_y} = \eta - \frac{v - (y_A - y_B)}{h_y}. \quad (25)$$

With (23)–(25), (22) is turned into

$$\begin{aligned} S_{1x}(u, v) &= S(\zeta, \eta) \\ &= \int_{D(\zeta)} \int_{D(\eta)} f(\zeta, \eta) g(\zeta - u, \eta - v) d\zeta d\eta \end{aligned} \quad (26)$$

with

$$\begin{aligned} -2h_x &\leq \zeta \leq 2h_x \\ -h_y &\leq \eta \leq h_y. \end{aligned} \quad (27)$$

Now, (26), in the case of xx interactions, can be reduced after some long calculations to the summation of several terms, basically of the following two kinds:

$$\begin{aligned} &2h_y \left(\int_{f(u)}^0 f'(\zeta) f''(\zeta) d\zeta \right) \\ &2h_y \left(\int_0^{g(u)} g'(\zeta) g''(\zeta) d\zeta \right) \end{aligned} \quad (28)$$

$f(u)$ and $g(u)$ have forms such as

$$1 \pm 2h_x \pm u \quad (29)$$

while $f'(\zeta)$, $f''(\zeta)$, $g'(\zeta)$, and $g''(\zeta)$ have forms such as

$$1 \pm \frac{\zeta}{2h_x}. \quad (30)$$

The solution of (28) with (29) and (30) leads to the following generical solution for (26):

$$S_{1x}(u, v) = (a_3u^3 + a_2u^2 + a_1u + a_0)(b_3v^3 + b_2v^2 + b_1v + b_0). \quad (31)$$

The best way to derive the coefficients a_i and b_i in (31) is to recall that (26) is a bidimensional convolution of two rooftop functions. If we indicate with z the distance between two interacting elementary cells, the solution of the convolution with graphical methods easily leads to the following formula:

$$S_{1x}(u, v) = h_x h_y R_{\zeta\zeta} \left(u - \frac{x_A - x_B}{h_x} \right) R_{\eta\eta} \left(v - \frac{x_A - x_B}{h_y} \right) \quad (32)$$

$R_{\zeta\zeta}(z)$ being the convolution along ζ , and $R_{\eta\eta}(z)$ being the convolution along η , with

$$R_{\zeta\zeta} = \begin{cases} \frac{z^3}{6} + z^2 + 2z + 4/3, & -2 \leq z \leq -1 \\ -\frac{z^3}{2} - z^2 + 2/3, & -1 \leq z \leq 0 \\ \frac{z^3}{2} - z^2 + 2/3, & 0 \leq z \leq 1 \\ -\frac{z^3}{6} + z^2 - 2z + 4/3, & 1 \leq z \leq 2 \\ 0, & \text{elsewhere} \end{cases} \quad (33)$$

$$R_{\eta\eta} = \begin{cases} z + 1, & -1 \leq z \leq 0 \\ 1 - z, & 0 \leq z \leq 1 \\ 0, & \text{elsewhere.} \end{cases} \quad (34)$$

Therefore, it is apparent that the coefficients a_i and b_i assume different values depending on the mutual distance z between the interacting cells, even though their form is analytically determined. For instance, using (9), for two cells with $z = 1$, we have

$$S_{1x}(r, \xi) = h_x h_y \left(1/2 \left(\frac{r \sin \xi - (x_A - x_B)}{h_x} \right)^3 - \left(\frac{r \sin \xi - (x_A - x_B)}{h_x} \right)^2 + 2/3 \right) \cdot \left(1 - \frac{r \cos \xi - (y_A - y_B)}{h_y} \right). \quad (35)$$

Once the S_{1x} has been determined, the W_{1x} is easily evaluated. In fact, we can now observe in (18) that S_{1x} is composed of terms $\cos^m \xi$ and $\sin^n \xi$, with $m, n = 0, 1, 2, 3$, whose primitives are known in closed form. For each value of r , $\xi_1(r)$ and $\xi_2(r)$ can be numerically evaluated: as apparent from Fig. 1, with simple calculations, the values for ξ can be found for all the interacting cells. Moreover, inside each cell the appropriate a_i and b_i can be identified with (32) to perform the integration. The domain of S_{1x} can be partitioned into eight subdomains [the four partitions of $R_{\zeta\zeta}$ in (33) and the two of $R_{\eta\eta}$ in (34)], and the calculation to evaluate W_{1x} become extremely fast.

In conclusion, $W_{1x}(r)$ is evaluated in a very efficient and nearly completely analytical way. Similar procedures can be followed to evaluate $W_{2x}(r)$. It can be derived that

$$S_{2x}(u, v) = \frac{h_y}{h_x} R_{\zeta\zeta} \left(u - \frac{x_A - x_B}{h_x} \right) R_{\eta\eta} \left(v - \frac{y_A - y_B}{h_y} \right). \quad (36)$$

$R_{\zeta\zeta}(z)$ being the convolution along ζ and $R_{\eta\eta}(z)$ being the convolution along η with

$$R_{\zeta\zeta} = \begin{cases} -z - 2, & -2 \leq z \leq -1 \\ 3z + 2, & -1 \leq z \leq 0 \\ -3z + 2, & 0 \leq z \leq 1 \\ z - 2, & 1 \leq z \leq 2 \\ 0, & \text{elsewhere} \end{cases} \quad (37)$$

$$R_{\eta\eta} = \begin{cases} z + 1, & -1 \leq z \leq 0 \\ 1 - z, & 0 \leq z \leq 1 \\ 0, & \text{elsewhere.} \end{cases} \quad (38)$$

The use of (9) transforms S_{2x} into a summation of terms $\cos^m \xi$ and $\sin^n \xi$ with $m, n = 0, 1, 2, 3$ (as in the previous case), whose primitives are known in closed form, thus making the derivation of W_{2x} quite immediate.

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