Scattering From Quasi-Planar and Moderate Rough Surfaces: Efficient Method to Fill the EFIE-Galerkin MoM Impedance Matrix and to Solve the Linear System

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Abstract—First, an acceleration to compute the impedance matrix obtained from the electric field integral equation (EFIE) discretized by the Galerkin method of moments (MoM) with Rao–Wilton–Glisson basis functions is addressed. It is based on a far-field approximation and makes it possible to avoid looping on the source and observation triangles. Next, the impedance matrix is split into strong and weak interactions; the latter is compressed by expressing it from Toeplitz submatrices. Then, the linear system is efficiently solved by a bi-iterative scheme. For a given order, the least-squares QR (LSQR) algorithm is applied to solve the sparse linear system related to the strong interactions, while the matrix–vector products, related to the weak interactions, are accelerated by using FFTs. Numerical results of the field scattered by perfectly conducting paraboloidshape object and Gaussian rough surface are shown.

Index Terms—Electric field integral equation (EFIE), fast algorithm, method of moments (MoM), radar cross section, rough surface scattering.

I. INTRODUCTION

THE calculation of full vector wave scattering from a large perfectly conducting surface is a very challenging issue; the main difficulty lies in the problem size. Solving a problem of this kind, from the electric field integral equation (EFIE) discretized by the Galerkin method of moments (MoM) with Rao–Wilton–Glisson basis functions [1], requires a great number of unknowns N_{Edge} in order to obtain accurate and meaningful results. Thus, reduction of both computation time and data storage requirement is continuously in progress.

Direct solvers, such as the lower upper (LU) decomposition, require $\mathcal{O}(N_{\text{Edge}}^3)$ operations while iterative solvers [2], such as conjugate gradient or generalized minimal residual (GMRES) techniques [3], need $\mathcal{O}(N_{\text{Edge}}^2)$ operations for the matrix–vector multiplication at each iteration. The memory requirement for these two solvers is usually $\mathcal{O}(N_{\text{Edge}}^2)$. Such computational complexity and memory requirements are too restrictive to be able to solve a large-scale scattering problem.

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These two disadvantages can be alleviated by using the sparse matrix canonical grid method (SMCG) [4]-[7], the adaptive integral method (AIM) [8]–[12], the stabilized extended boundary condition method (SEBCM) [13], and the fast multipole method (FMM) [14], [15]. For far-field interactions, the principle of AIM and SMCG is similar and consists of expressing the Green function on a uniform grid by using an interpolation scheme. From SMCG, a Taylor series expansion is also applied on the surface elevations. It is related to the concept of the short interaction range [16]. This implies that the matrix is Toeptliz, making it possible to reduce the memory requirement and to accelerate the matrix-vector product by using FFTs. The first novelty of this article is to accelerate the calculation of the weak impedance matrix elements by deriving a closed-form expression, resulting from the sum over the four facets common to the source and observation edges. In addition, since with a triangular discretization the basis functions are not uniformly distributed, the weak impedance matrix is expressed from three Toeplitz submatrices by defining two uniform subgrids. This avoids applying interpolations. Next, for quasi-planar surfaces and similar to [7], the surface elevation Δz is accounted for by expanding the Toeplitz submatrices over Δz .

For an iterative solver, the convergence of iterations is not very good for ill-posed matrix equations. This often occurs when the EFIE is used to solve 3-D complex scattering problems. For instance, Hu and Nie [17] added the MFIE contribution to decrease the condition number of the resulting matrix, and the linear system is solved from a bi-iterative procedure. The present method distinguishes the strong nearfield interactions and the weak ones between the observation point and the source point. This has the advantage that only the near interaction matrix is stored, and the far interaction matrix is Toeplitz by blocks and requires to store only $\mathcal{O}(N_{\text{Edge}})$ elements. The second novelty of this article is to take advantage of this decomposition for solving efficiently the linear system from a bi-iterative algorithm. Then, for a given order, the least-squares QR (LSQR) algorithm [18] is applied to efficiently solve the sparse linear system related to the strong interactions, while the matrix-vector products, related to the weak interactions, are accelerated by using FFTs. Numerical tests showed that LSQR converges

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more rapidly than GMRES or biconjugate gradient stabilized method.

This article is organized as follows. Section II presents the EFIE impedance matrix and the scattered field. Section III addresses the derivation of the weak interaction impedance matrix. Section IV deals with how this matrix can be partitioned into Toeplitz submatrices. Section V presents the bi-iterative scheme to efficiently solve the linear system. Section VI shows numerical results for a paraboloid-shaped object and a Gaussian rough surface. Section VII gives concluding remarks.

II. EFIE IMPEDANCE MATRIX AND SCATTERED FIELD

In this article, to compute the field scattered by a perfectly conducting object, the EFIE is solved from the MoM. In addition, the Galerkin method is applied by using the Rao–Wilton–Glisson basis functions. This leads to solving the linear system $\bar{Z}X = b$, where \bar{Z} is the impedance matrix and X a vector related to the incident wave. The time convention $e^{-j\omega t}$ is used throughout this article.

The element $Z_{m,n}$ of the impedance matrix $\overline{\mathbf{Z}}$, corresponding to the interaction between two edges *m* (observation) and *n* (source) of a facet couple (p, q), is expressed as [2]

$$Z_{m,n} = \frac{s_p s_q}{4\pi A_p A_q L_m L_n} \iint_{T_p} \iint_{T_q} \left(\rho_m^p \cdot \rho_n^q - \frac{1}{k^2} \right) \\ \times \frac{e^{-jkD_{p,q}}}{D_{p,q}} dR_p dR_q \quad (1)$$

where $s_{p,q} = \pm 1$ and $\{A_{p,q}\}$ are the triangle areas, $\{L_{m,n}\}$ the edge lengths, and $\rho_{m,n}^{p,q} = (V_{m,n}^{p,q} - R_{p,q})/2$, in which $V_{m,n}^{p,q}$ is the position vector of the vertex unshared by the edge (m, n) and belonging to the facet (p, q). In addition, $D_{p,q} = ||\mathbf{R}_p - \mathbf{R}_q||$ and k is the wavenumber that equals $2\pi/\lambda$, where λ is the wavelength in free space.

Assuming a plane incident wave, a component b_m of the vector **b** associated with the source edge *m* and facet *p* is given by [2]

$$b_m = -\frac{j}{\omega\mu} \frac{L_m s_m}{2A_p} \iint_{\mathbf{T}_p} \boldsymbol{\rho}_m^p \cdot \hat{\boldsymbol{p}}_{\mathrm{inc}} e^{-\boldsymbol{k}_{\mathrm{inc}} \cdot \boldsymbol{R}_p} dR_p \qquad (2)$$

where ω is the wave pulsation and μ the permeability of the surrounding medium. In addition, \hat{p}_{inc} [either vertical, \hat{v}_{inc} (θ), or horizontal, \hat{h}_{inc} (ϕ)] and k_{inc} are the polarization and incident wave vectors, respectively, both defined in spherical coordinates from the angles (θ_{inc} , ϕ_{inc}). By solving the linear system $X = \bar{Z}^{-1}b$, the components { a_n } of the vector X are found. The scattered far-field is then expressed as

$$\boldsymbol{E}_{\mathrm{sca}}^{\infty}(\boldsymbol{R}_{0}) = -\frac{j\omega\mu e^{-j\boldsymbol{k}\boldsymbol{R}_{0}}}{8\pi\,\boldsymbol{R}_{0}}\sum_{p=1}^{P_{\mathrm{Fact}}}\sum_{m=1}^{M_{\mathrm{Edge}}}\frac{L_{m}a_{m}s_{m}}{A_{p}}$$
$$\times \iint_{\mathbf{T}_{p}}\boldsymbol{\rho}_{m}^{p}e^{j\boldsymbol{k}_{\mathrm{sca}}\cdot\boldsymbol{R}_{p}}d\boldsymbol{R}_{p} \quad (3)$$

where P_{Facet} is the number of facets and M_{Edge} the number of edges associated with the facet *p*. In addition, R_0 is the

distance from the receiver to the phase origin of the object. The scattering coefficient is then expressed as

$$SC_{p_{inc}p_{sca}} = \lim_{R_0 \to \infty} 2\sqrt{\pi} R_0 \frac{E_{sca}^{\infty} \cdot \hat{p}_{sca}}{E_{inc} \cdot \hat{p}_{inc}}$$
(4)

where $p_{inc} = \{\theta, \phi\}$ and $p_{sca} = \{\theta, \phi\}$. The subscripts "inc" and "sca" stand for incident and scattered (waves), respectively. The receiver polarization basis ($\hat{k}_{sca}, \hat{v}_{sca}, \hat{h}_{sca}$) can be defined in a similar way as that of the incident field ($\hat{k}_{inc}, \hat{v}_{inc}, \hat{h}_{inc}$), in which θ_{sca} and ϕ_{sca} are the receiver (scattering) angles. The radar cross section RCS_{*p*inc *p*sca} is obtained by taking the squared modulus of SC_{*p*inc *p*sca}.

III. DERIVATION OF THE WEAK INTERACTION IMPEDANCE MATRIX

The impedance matrix is split into near \bar{Z}_{Strong} (or strong) and far \bar{Z}_{Weak} (or weak) interactions as

$$\mathbf{Z} = \mathbf{Z}_{\text{Strong}} + \mathbf{Z}_{\text{Weak}} \tag{5}$$

where the elements of the strong interactions are calculated from (1) (without approximation) and those of the weak interactions from the equation derived in Appendix A. Equation (1) shows that $Z_{m,n}$ requires the calculation of two twofold numerical integrations over the surfaces of the triangles T_p and T_q . This is done from twofold Gauss–Legendre integrations. In this article, the weak interactions are derived from a closedform expression addressed in Appendix A. The singularity, which occurs for $D_{p,q} = 0$, is computed from the work published by [19].

Then, calculating the sum over the four facets of edges (m, n), Appendix A shows that an element of the weak interaction is

$$\tilde{Z}_{m,n} \approx \sum_{p=1}^{p=3} G^{(p)}(\boldsymbol{R}_m - \boldsymbol{R}_n) W^{(p)}(\boldsymbol{R}_m, \boldsymbol{R}_n) = \sum_{p=1}^{p=3} G^{(p),(0)}(r_{m,n}) \sum_{q=0}^{q=Q} W^{(p)}(\boldsymbol{R}_m, \boldsymbol{R}_n) \times (z_m - z_n)^{2q} A^{(p),(q)}(r_{m,n})$$
(6)

where $\mathbf{R}_{m,n} = \mathbf{r}_{m,n} + (z_m - z_n)\hat{z}$ and $G_{m,n}^{(p),(0)} = G_{m,n}^{(p)}|_{z_n-z_m=0}$. The above equation shows that the matrix $\tilde{\mathbf{Z}}_{m,n}$ is expressed as the sum over three matrices { $\tilde{\mathbf{G}}^{(p)}$ } that depend only on $\mathbf{R}_m - \mathbf{R}_n$ weighted by polynomial functions { $W^{(p)}$ } that depend on the source edge \mathbf{R}_n (its middle) and observation edge \mathbf{R}_m (its middle). In addition, the element $G_{m,n}^{(p)}$ can be expanded over $z = z_m - z_n$ near 0, where z is the elevation difference between the edge middles m and n. It is important to underline that $G_{m,n}^{(p),(0)}$ and $A_{m,n}^{(p),(q)}$ (weighting of the Taylor series expansion) depend only on $\mathbf{r}_{m,n} = \sqrt{(x_m - x_n)^2 + (y_m - y_n)^2}$. The functions $W_{m,n}^{(p)}$, $G_{m,n}^{(p),(0)}$, and $A_{m,n}^{(p),(q)}$ are expressed from (A15), (A17), (A20) (Q = 1) and (A21) (Q = 2).

The appendix also shows that (6) is valid if the strong interaction distance ((A6) and (A22)) satisfies

$$r_{\text{Strong}} > \max\left(\frac{\max(L_m, \Delta z)^2 n_0}{2\lambda}, \max(\Delta z) \sqrt{\frac{n_0}{2}}\right)$$
 (7)

where $n_0 = 20$, $\Delta z = \max(z) - \min(z)$ on the distance r_{Strong} , and L_m is the *m* edge length.

If the (x_m, y_m) (or (x_n, y_n)) Cartesian grid is uniform, then the matrix associated with the element $G_{m,n}^{(p),(0)} \times A_{m,n}^{(p),(q)} = B_{m,n}^{(p),(q)}$ is Toeplitz. In addition, since $W_{m,n}^{(p)} = \sum_{s=1}^{s=S} f_m^{(s)} g_n^{(s)}$ is a polynomial function of \mathbf{R}_n and \mathbf{R}_m , (6) can be written as

$$\tilde{Z}_{m,n} \approx \sum_{p=1}^{p=3} \sum_{q=0}^{q=Q} \sum_{r=0}^{R=2q} W_{m,n}^{(p)} B_{m,n}^{(p),(q)} (z_m)^{2q-r} (-z_n)^r C_{2q}^r$$

$$= \sum_{p=1}^{p=3} \sum_{s=1}^{s=S} \sum_{q=0}^{q=Q} \sum_{r=0}^{R=2q} C_{2q}^r [(z_m)^{2q-r} f_m^{(s)}] B_{m,n}^{(p),(q)}$$

$$\times [(-z_n)^r g_n^{(s)}]$$
(8)

where $B_{m,n}^{(p),(q)} = G^{(p),(0)}(r_{m,n})A^{(p),(q)}(r_{m,n})$ and $C_{2q}^r = (2q)!/[r!(2q-r)!]$ (binomial coefficient). The above equation is in a form such that the source point *n* is on the righthand side of $B_{m,n}^{(p),(q)}$, while the observation point *m* is located on the left-hand side of $B_{m,n}^{(p),(q)}$. Appendix A shows that S = 3 + 6 + 36 = 45 and gives the expression of $\{g_n^{(s)}, f_m^{(s)}\}$ for $s = \{1, 2, 3\}$, corresponding to p = 1. The terms defined for s > 3 are obtained in a similar way.

The integer *S* corresponds to the expansion order that decomposes any function as a sum of *S* terms that depend only on the (m, n) product. For example, the dot product $\rho_m \cdot \rho_n = \rho_{m,x}\rho_{n,x} + \rho_{m,y}\rho_{n,y} + \rho_{m,z}\rho_{n,z}$, where the subscripts (x, y, z) denote the vector components. Then S = 3. This way is applied in (A15).

Then, the matrix product $O\overline{Z}S$ can be computed from $3 \times 45 \times [Q(Q+2)+1]$ FFTs, where O and S are any observation and source vectors, respectively.

IV. TOEPLITZ SUBMATRICES

As shown in Fig. 1, the (x_m, y_m) (or (x_n, y_n)) Cartesian grid over the center of the edges is not uniform, but the grid can be uniform by blocks. To this end, as shown in Fig. 1, the edges are sorted in ascending order going from left to right and from bottom to top. In what follows, the edges of y values equal to $\{-0.15, -0.5, 0.5, 0.15\}\lambda$ are named *odd* rows, whereas those of y values equal to $\{-0.1, 0, 0.1\}\lambda$ are named *even* rows.

Let *f* be any function of $r_{m,n} = \sqrt{(x_n - x_m) + (y_n - y_m)^2}$ (planar surface) and $Z_{m,n}$ an element of the matrix $\bar{\mathbf{Z}}$ associated with the function $f(r_{m,n})$. Then, as shown in Fig. 2, $\bar{\mathbf{Z}}$ is Toeplitz by blocks by considering four cases.

- 1) The purple color represents the Toeplitz submatrices between the edges of *odd* rows.
- 2) The blue color represents the Toeplitz submatrices between the edges of *even* rows.
- 3) The green color represents the Toeplitz submatrices between the edges of odd rows and even rows, for which the edge numbers are *odd*.
- 4) The yellow color represents the Toeplitz submatrices between the edges of odd rows and even rows, for which the edge numbers are *even*.

In addition, for two rows separated by the same y distance, the matrix is the same. For example, in Fig. 2, the submatrix of edges $m = n = \{1, 2, 3, 4, 5, 6, 7\}$ is the same as



Fig. 1. Sort of the edges. A square plate of area $L_x \times L_y = (0.4\lambda)^2$ is considered, and the sampling steps with respect to the *x*- and *y*-directions are $\Delta x = \Delta y = 0.1\lambda$. The label $m \Rightarrow m'$ (or $n \Rightarrow n'$) means that the first number *m* indicates the original edge number and the second number *m'* gives the new edge number. The integer *m'* sorts the edges in ascending order going from left to right and from bottom to top.



Fig. 2. Structure of the impedance matrix represented by Toeplitz submatrices. The four cases are represented by four colors, and the geometry is shown in Fig. 1.

that obtained for $m = n = \{12, 13, 14, 15, 16, 17, 18\}$. For cases 1, 2, and 3, the submatrices are also symmetric, which implies that only the first row is calculated, whereas, for case 4, the first row and column are computed.

For each case, Table I lists the length of the rows or columns of the submatrices and their number. $N_{x,y}$ is the number of vertices on a uniform grid of step Δx and Δ_y , respectively. For instance, in Fig. 1, $\Delta x = \Delta y = 0.1\lambda$, $N_x = N_y = 5$.

The total number of elements is then $N_{\text{Toep}} = 6N_x$ $N_y - 10N_x - 8N_y + 13$, while the number of edges is

Case	Number	Length
1	$N_y - 1$	$2N_x - 3$
2	$N_y - 2$	$N_x - 1$
3	$N_y - 2$	$N_x - 1$
4	$N_{y} - 2$	$N_x - 1$
4	$N_y - 2$	$N_x - 2$

 $N_{\text{Edge}} = 3N_x N_y - 4(N_x + N_y) + 5$. For $N_x \gg 1$ and $N_y \gg 1$, $N_{\text{Toep}} \approx 2N_{\text{Edge}}$. For the weak interactions and a planar surface, this means that, instead of calculating $N_{\text{Edge}}^2 - N_{\text{Strong}}$ elements, only $2N_{\text{Edge}}$ elements are computed, where N_{Strong} is the number of edge pairs in near field. In other words, for $N_{\text{Edge}}^2 \gg N_{\text{Strong}}$, the compression rate is of the order of $1 - 2/N_{\text{Edge}}$.

In addition, from the Section III and Table I, for $N_{x,y} \gg 1$, the complexity of the matrix-vector product $\bar{Z}X = (\bar{Z}_{\text{Strong}} + \bar{Z}_{\text{Weak}})X$ is

$$C_{\text{MVP}} = 135 \times [Q(Q+2)+1] \times 4 \times N_y \log_2 N_x + N_{\text{Strong}}$$

= 540[Q(Q+2)+1]N_y log_2 N_x + N_{\text{Strong}}
$$\approx \alpha \sqrt{N_{\text{Edge}}} \log_2 N_{\text{Edge}} + N_{\text{Strong}}$$
(9)

where $\alpha = 270[Q(Q+2)+1]/\sqrt{3}$, $N_{\text{Edge}} \approx 3N_x N_y = 3N_x^2$ with $N_y = N_x$. For $r_{\text{Strong}} = 0$, $N_{\text{Strong}} = N_{\text{Edge}}^2$ ($\bar{\mathbf{Z}}_{\text{Weak}} = \bar{\mathbf{0}}$). Then, in comparison to a conventional matrix-vector product of complexity N_{Edge}^2 , the use of FFTs is efficient if

$$\eta_{\rm MVP} = \frac{\alpha \log_2 N_{\rm Edge}}{N_{\rm Edge}^{3/2}} + \frac{N_{\rm Strong}}{N_{\rm Edge}^2} \ll 1. \tag{10}$$

V. RESOLUTION OF THE LINEAR SYSTEM

The final step is to efficiently solve the linear system $\bar{Z}X = b$. For large problems, the conventional LU decomposition cannot be used, and iterative schemes are preferred. Usually, the conjugate gradient algorithm and their improved versions can be good candidates, but, for the EFIE, their convergence order is very large (for the scenarios presented in Section VI, the order exceeds 300). In this article, we propose to use \bar{Z}_{Strong} as a preconditioning matrix $\bar{M}_c = \bar{Z}_{\text{Strong}}^{-1} \bar{Z}_{\text{Weak}}$, and next, a Taylor series expansion is applied on the resulting characteristics matrix to account for the weak interactions through \bar{Z}_{Weak} . Then

$$X = (\bar{Z}_{\text{Strong}} + \bar{Z}_{\text{Weak}})^{-1} b$$

= $\left[\bar{Z}_{\text{Strong}} \left(\bar{I} + \bar{Z}_{\text{Strong}}^{-1} \bar{Z}_{\text{Weak}}\right)\right]^{-1} b$
= $(\bar{I} + \bar{M}_c)^{-1} \bar{Z}_{\text{Strong}}^{-1} b$
 $\approx \sum_{k=0}^{K} (-\bar{M}_c)^k \bar{Z}_{\text{Strong}}^{-1} b = \sum_{k=0}^{K} \bar{Y}^{(k)}$ (11)

in which

$$\begin{cases} \bar{\boldsymbol{Y}}^{(0)} = \bar{\boldsymbol{Z}}_{\text{Strong}}^{-1} \boldsymbol{b} & k = 0\\ \bar{\boldsymbol{Y}}^{(k+1)} = -\bar{\boldsymbol{M}}_{c} \bar{\boldsymbol{Y}}^{(k)} = -\bar{\boldsymbol{Z}}_{\text{Strong}}^{-1} \left(\bar{\boldsymbol{Z}}_{\text{Weak}} \bar{\boldsymbol{Y}}^{(k)} \right) & k > 0 \end{cases}$$
(12)



Fig. 3. Paraboloid surface with $L_x = L_y = 12\lambda$. $N_{\text{Edge}} = 42,960, \Delta x = \Delta y = 0.1\lambda$, and $a = \lambda/2$.

This algorithm converges if the spectral radius (largest modulus of its eigenvalues) of \bar{M}_c is strictly smaller than one. This implies that the strong distance must be not too small. For large problems, this value is not calculated because it is very time-consuming. Since \bar{Z}_{Strong} is a sparse matrix, its LU decomposition matrices are also sparse, but this operation can be time-consuming. To overcome this issue, the LSQR [18] (least-squares QR) algorithm is applied because it is efficient for a sparse matrix and more efficient than a gradient conjugate-based method.

From (12), the complexity of the I-LSQR (I as iterative) algorithm is

$$C_{\text{I-LSQR}} = C_{\text{LSQR}} + K_{\text{I-LSQR}} (C_{\text{LSQR}} + C_{\text{MVP}})$$
(13)

where C_{LSQR} is the complexity of the LSQR algorithm and C_{MVP} is expressed from (9). The convergence order $K_{I-LSQR} = k$ is obtained when the relative residual error (RRE) satisfies

$$\operatorname{RRE} = \frac{\operatorname{norm}(X^{(k+1)} - X^{(k)})}{\operatorname{norm}(X^{(k+1)})} < \epsilon_{\operatorname{I-LSQR}}$$
(14)

where $X^{(K)} = \sum_{k=0}^{k=K} Y^{(k)}$ and $\epsilon_{\text{I-LSQR}}$ is the threshold of I-LSQR. Typically, $\epsilon_{\text{I-LSQR}} = 10^{-2}$ and $\epsilon_{\text{LSQR}} = 5 \times 10^{-4}$ equal the LSQR threshold.

VI. NUMERICAL RESULTS

The wavelength in free space λ is 1 m, and the polarization is $\theta\theta$.

A. Paraboloid Surface

First, a paraboloid surface of equation z(x, y) = 2a $(x^2 + y^2)/L_x^2$ is considered, and it is shown in Fig. 3. In the (Ox, Oy) plane, the surface area is $L_x L_y$, where $\{L_{x,y}\}$ are the surface lengths with respect to the x- and y-directions. In addition, the center of the surface is the point O of coordinates (0, 0). For $x = L_x/2$ and $y = L_y/2 = L_x/2$, $z = a = \max(z)$. In Fig. 3, the number of edges is $N_{\text{Edge}} = 42,960$, and the sampling steps with respect to the x- and y-directions are $\Delta x = \Delta y = 0.1\lambda$ and $a = \lambda/2$.

Fig 4 shows the plots of the bistatic RCS in dBm² versus the scattering angle θ_{sca} . To better highlight the differences,



Fig. 4. Top: bistatic RCS in dBm² versus the scattering angle θ_{sca} . Bottom: ratio RCS_{1-LSQR}/RCS_{LU} in dB scale. Q = 0, $r_{Strong} = \{1.2, 1.4, 1.6\}\lambda$, $\theta_{inc} = \pi/6$, $\phi_{inc} = 0$, and $\phi_{sca} = 0$. The illuminated object is a paraboloid surface, as shown in Fig. 3.

in the lower subfigure, the ratio RCS_{I-LSQR}/RCS_{LU} is plotted in dB scale, where RCS_{LU} is the RCS computed from a LU decomposition of the impedance matrix. The incidence angles are $\theta_{inc} = \pi/6$ and $\phi_{inc} = 0$, and $\phi_{sca} = 0$ is the azimuthal scattering angle. The illuminated object is a paraboloid surface, as shown in Fig. 3. In the legend, "LU: $(t_{LU,1}, t_{LU,2})$ s" and " u_{Strong} , I-LSQR – $(Q, K_{I-LSQR}, N_{LSQR})$: $(t_{I-LSQR,1}, t_{I-LSQR,2}, t_{I-LSQR,3})$ s, RRE" mean the following.

- 1) $t_{LU,1}$: Computing time to fill the impedance matrix, in seconds.
- t_{LU,2}: Computing time to solve the linear system from LU, in seconds.
- 3) $u_{\text{Strong}} = r_{\text{Strong}}/\lambda$: Distance of the strong interactions normalized by λ .
- 4) K_{I-LSQR} : Convergence order of I-LSQR.
- 5) N_{LSQR} : Mean convergence order of LSQR ($\epsilon_{\text{LSQR}} = 5 \times 10^{-4}$).
- 6) Q: Order of the Taylor series expansion over z.
- 7) $t_{I-LSQR,1}$: Computing time to fill the impedance matrix of the strong interactions, in seconds.
- 8) $t_{\text{I-LSQR},2}$: Computing time to compute the matrix-vector product $\bar{Z}_{\text{Weak}}\bar{Y}^{(k)} = v$, in seconds.
- 9) $t_{\text{I-LSQR},3}$: Computing time to solve the linear system $\bar{Z}_{\text{Strong}}^{-1} v$ from LSQR ($\epsilon_{\text{LSQR}} = 5 \times 10^{-4}$), in seconds.
- 10) RRE: RRE obtained at the convergence order $K_{\text{I-LSQR}}$ ($\epsilon_{\text{I-LSQR}} = 10^{-2}$).

The total computing time is then $t_{LU,1} + t_{LU,2}$ for LU and $t_{I-LSQR,1} + t_{I-LSQR,2} + t_{I-LSQR,3}$ for I-LSQR.

As we can see in Fig. 4, the results match well with those obtained from LU. As r_{Strong} increases, the results better match, the order of convergence, $K_{\text{I-LSQR}}$, decreases from 5 to 3, and the filling time $t_{\text{I-LSQR},1}$ increases slightly. By calculating only \bar{Z}_{Strong} , the memory requirement is divided by 20 in comparison to LU, which needs to calculate all the elements of \bar{Z} . Compared to the filling computation time of LU, the gain is



Fig. 5. Same variations as in Fig. 4, but $a = 1.5\lambda$, $r_{\text{Strong}} = 1.6\lambda$, and $Q = \{0, 1, 2\}$.

of the order of 9. On the other hand, the time $t_{I-LSQR,2}+t_{I-LSQR,3}$ to solve the linear system by I-LSQR is slightly smaller than that of LU, $t_{LU,2}$.

From (10), $\eta_{\text{MVP}} \approx 0.05$, with $r_{\text{Strong}} = 1.6\lambda$. This means that the matrix–vector product computed from N_{FFT} FFTs should be $1/\eta_{\text{MVP}} \approx 20$ faster than the one computed in a conventional manner. In practice, it is not the case. The FFTs [and inverse fast Fourier transforms (IFFTs)] are computed from the fft and ifft MATLAB functions in a matrix manner to accelerate their calculation. Nevertheless, loops are required, and the reshape MATLAB function is also applied, which increases the computing time. If this step was programed in C, the computing time would be smaller. We can also note that this time $t_{\text{I-LSQR},2}$ is larger than $t_{\text{I-LSQR},3}$, allocated to solve the linear system by LSQR, also computed in MATLAB but by calling C functions.

However, both computing times ($t_{I-LSQR,2}$, $t_{I-LSQR,3}$) are not comparable, and $t_{I-LSQR,2}$ would be much smaller if the matrix–vector product was made in C. In addition, a parallelization would make it possible to significantly to decrease this computation time. With MATLAB, the LU inversion is optimized and parallelized.

Fig. 5 shows the plots of the same variations as in Fig. 4, but $a = 1.5\lambda$, $r_{\text{Strong}} = 1.6\lambda$, and $Q = \{0, 1, 2\}$. As expected, as Q increases, the results better match with those obtained from LU, and the results computed for Q = 1 and Q = 2 are nearly the same.

Equation (7) gives the lowest value of the strong interaction distance, for which the derivation of the weak interaction matrix is valid at order zero (Q = 0). In polar coordinates, $z(x, y) = h(r) = 2ar^2/L_x^2$ and $dh/dr = 4ar/L_x^2$. Then, $\max(\Delta z) \approx 4a \max(r)r_{\text{Strong}}/L_x^2 = 2ar_{\text{Strong}}/L_x = 0.4\lambda$ in Fig. 5. From (7), $r_{\text{Strong}} > \max(0.22, 1.60, 126)\lambda = 1.6\lambda =$ $r_{\text{Strong,min}}$ with $n_0 = 20$. This explains in Fig. 5 why the results for Q = 0 deviate from those obtained from LU for low values of the RCS since $r_{\text{Strong}} = r_{\text{Strong,min}}$. The order Q = 1makes it possible to significantly decrease this deviation, and the contribution of the second order is negligible.



Fig. 6. Surface height versus the coordinates x and y. $L_x \times L_y = 144\lambda^2$, $\sigma_z = 0.3\lambda$, and $L_{c,x} = L_{c,y} = 1.5\lambda$. The number of edges is $N_{\text{Edge}} = 42\,960$.



Fig. 7. Top: bistatic RCS in dBm² versus the scattering angle θ_{sca} . Bottom: ratio RCS_{1-LSQR}/RCS_{LU} in dB scale. Q = 0, $r_{Strong} = \{1.2, 1.4, 1.6\}\lambda$, $\theta_{inc} = \pi/6$, $\phi_{inc} = 0$, and $\phi_{sca} = 0$. The illuminated object is a rough surface, which is plotted in Fig. 6.

B. Rough Surface

To strongly attenuate the edge diffractions by the surface, the well-known incident tapered wave published by Braunish *et al.* [20] is applied (at the order two) with tapering parameter $g = L_x/4$ (surface of area $A_0 = L_x^2$). The length *g* controls the extent of the incident beam that illuminates the surface.

Fig. 6 shows the plots of a random rough surface of Gaussian height distribution and Gaussian height autocorrelation function. The surface correlation lengths with respect to the *x*- and *y*-directions are $L_{c,x} = L_{c,y} = 1.5\lambda$, the surface height standard deviation is $\sigma_z = 0.3\lambda$, and the surface area is $L_x \times L_y = 144\lambda^2$.

Fig. 7 shows the plots of the bistatic RCS in dBm² versus the scattering angle θ_{sca} . At the bottom, the ratio RCS_{I-LSQR}/RCS_{LU} is plotted in dB scale. Q = 0, $r_{Strong} = \{1.2, 1.4, 1.6\}\lambda$, $\theta_{inc} = \pi/6$, $\phi_{inc} = 0$, and $\phi_{sca} = 0$. The illuminated object is a rough surface, which is plotted in Fig. 6. For high values of RCS, the results match well with those obtained from LU. From (7) and by taking $\Delta z \approx \sqrt{2}\sigma_z$,



Fig. 8. Same variations as in Fig. 7, but the incidence angle is $\theta_{inc} = \pi/4$.



Fig. 9. Same variations as in Fig. 7, but $r_{\text{Strong}} = 1.6\lambda$ and $Q = \{0, 1, 2\}$.

 $r_{\text{Strong}} > \max(0.34, 1.80, 1.34)\lambda = 1.8\lambda$. In Fig. 7, this explains why differences with LU appear for low values of RCS.

Fig. 8 shows the plots of the same variations as in Fig. 7, but the incidence angle is $\theta_{inc} = \pi/4$ (instead of $\theta_{inc} = \pi/6$). Like in Fig. 7, a deviation occurs from incidence angles larger than 30°–45°, and the proposed method gives similar performances.

Fig. 9 shows the plots of the same variations as in Fig. 7, but $r_{\text{Strong}} = 1.6\lambda$ and $Q = \{0, 1, 2\}$. For Q = 1, the results match well with those obtained from LU. Fig. 9 also shows that it is not relevant to calculate the order Q = 2.

Fig. 10 shows the plots of the same variations as in Fig. 9 but in cross polarization $[\theta\phi]$ or vertical-horizontal (VH)]. The VH strengths are 20 dB lower than those obtained in vertical-vertical (VV), which explains why the deviation with LU is larger than those shown in Fig. 9. Indeed, the cross polarization requires better accuracy since the levels are smaller.

Fig. 11 shows the plots of the same variations as in Fig. 9, but $L_{c,x} = L_{c,y} = 2\lambda$ (instead of $L_{c,x} = L_{c,y} = 1.5\lambda$). As we can see, the results obtained for Q = 1 match well



Fig. 10. Same variations as in Fig. 9 but in cross polarization ($\theta\phi$ or VH).



Fig. 11. Same variations as in Fig. 9, but $L_{c,x} = L_{c,y} = 2\lambda$ (instead of $L_{c,x} = L_{c,y} = 1.5\lambda$).



Fig. 12. Same variations as in Fig. 9, but $r_{\text{Strong}} = 4.4\lambda$ and $\sigma_z = 0.5\lambda$ (instead of $\sigma_z = 0.3\lambda$).

with those obtained from LU. In comparison to Fig. 9, the total computing time is slightly smaller because r_{Strong} remains

unchanged and the mean convergence order $N_{I-LSQR} \approx 117$ is smaller, whereas the convergence order $K_{LSQR} = 4$ remains constant.

Fig. 12 shows the plots of the same variations as in Fig. 9, but $r_{\text{Strong}} = 4.4\lambda$ and $\sigma_z = 0.5\lambda$ (instead of $\sigma_z = 0.3\lambda$). The value $r_{\text{Strong}} = 4.4\lambda$ is chosen so that it satisfies criterion (7). As we can see, the results obtained for Q = 1 match well with those computed from LU. In comparison to LU (and Fig. 9), the computing time is greater because r_{Strong} is larger than that used for $\sigma_z = 0.3\lambda$ (it is proportional to σ_z^2) and the mean convergence order $N_{I-\text{LSQR}} \approx 206$ increases, whereas $K_{\text{LSQR}} = 3$ is smaller. As expected, as σ_z increases, the proposed method is less efficient.

VII. CONCLUSION

First, an acceleration to compute the impedance matrix, based on a far-field approximation, is addressed. Next, the impedance matrix is split into strong and weak interactions, and this latter is compressed by expressing it from Toeplitz submatrices. Then, the linear system is efficiently solved from a bi-iterative scheme. For a given order, the LSQR algorithm is applied to solve the sparse linear system related to the strong interactions, while the matrix–vector products, related to the weak interactions, are accelerated by using FFTs.

For a paraboloid-shaped object, the numerical results show that the proposed method, named I-LSQR, is very efficient. The computation of the higher order (related to $\Delta z = \max(z) - \min(z)$) makes it possible to obtain more accurate results with an increase in the computing time.

For a rough surface, the numerical results also show that I-LSQR is efficient, but it is more sensitive to $\Delta z \propto \sigma_z$. In addition, the method becomes less efficient as σ_z increases because the minimum strong interaction distance (7) increases. It is approximately expressed as $(\sqrt{2}\sigma_z)^2 \times 20/(2\lambda) = 20\sigma_z^2/\lambda$.

APPENDIX A Derivation of the Weak Impedance Matrix Elements

For a pair of facets in far-field from each other, using an updated Fraunhofer criterion, Bourlier [21], [22] showed that (1) can be simplified as

$$Z_{m,n} \approx \frac{s_p s_q e^{-jkR_{p,q}}}{4\pi A_p A_q R_{p,q} L_m L_n} \iint_{T_p} \iint_{T_q} \left(\boldsymbol{\rho}_m^p \cdot \boldsymbol{\rho}_n^q - \frac{1}{k^2} \right) \\ \times e^{-jk\hat{\boldsymbol{R}}_{p,q} \cdot \left(\delta_p - \delta_q \right)} dR_p dR_q \tag{A1}$$

where $\mathbf{R}_{p,q} = \mathbf{G}_p - \mathbf{G}_q$ and $\delta_{p,q} = \mathbf{M}_{p,q} - \mathbf{G}_{p,q}$, in which $\mathbf{G}_{p,q}$ is the gravity center of the facets p and q, respectively, and $\mathbf{M}_{p,q}$ the integration point on the facets p and q, respectively $(\mathbf{D}_{p,q} = \mathbf{M}_p - \mathbf{M}_q = \mathbf{R}_{p,q} + \delta_p - \delta_q)$. In addition, $\hat{\mathbf{R}}_{p,q} = \mathbf{R}_{p,q}/\mathbf{R}_{p,q}$, where $\mathbf{R}_{p,q} = \|\mathbf{R}_{p,q}\|$.

For the weak interactions, the integral can be evaluated from a single point chosen at the middle of the egde *m* or *n*, $P_{m,n}$. Equation (A1) becomes

$$Z_{m,n} \approx \frac{s_p s_q e^{-jkR'_{m,n}}}{4\pi R'_{m,n}L_m L_n} \left(\boldsymbol{\rho}_m^p \cdot \boldsymbol{\rho}_n^q - \frac{1}{k^2} \right)$$
(A2)



Fig. 13. For a planar surface, interaction of an edge pair (m, n), and their two facets. The facets are assumed to be identical.

where

$$R'_{m,n} = R_{m,n} \left[1 + \frac{(\boldsymbol{\delta}_m - \boldsymbol{\delta}_n) \cdot \boldsymbol{R}_{m,n}}{R_{m,n}^2} \right]$$
(A3)

 $\delta_{p,q} = \delta_{m,n} = P_{m,n} - G_{p,q}$, and $R_{p,q} = P_m - P_n = R_{m,n}$. In comparison to (A2) at the denominator, $R_{p,q}$ is changed by $R'_{p,q}$. Since $|(\delta_m - \delta_n) \cdot R_{m,n}|/R^2_{m,n} \ll 1$, and by using the approximation $1/(1 + x) \approx 1 - x \approx e^{-x}$ for $|x| \ll 1$, (A2) can be expressed as

$$Z_{m,n} \approx \frac{s_p s_q e^{-jkR_{m,n}}}{4\pi R_{m,n} L_m L_n} \left(\boldsymbol{\rho}_m^p \cdot \boldsymbol{\rho}_n^q - \frac{1}{k^2} \right) e^{j(\phi_1 + \phi_2)}$$
(A4)

where

$$\begin{cases} \phi_1 = -k\delta_m \hat{\boldsymbol{\delta}}_m \cdot \hat{\boldsymbol{R}}_{m,n} \left(1 - \frac{j}{kR_{m,n}} \right) \\ \phi_2 = +k\delta_n \hat{\boldsymbol{\delta}}_n \cdot \hat{\boldsymbol{R}}_{m,n} \left(1 - \frac{j}{kR_{m,n}} \right) \end{cases}$$
(A5)

Equation (A4) is valid if $\Delta^2/2R_{m,n} < \lambda/n_0$ [21], where n_0 is an integer ranging from 10 to 20 and $\Delta = \max(\delta_n) + \max(\delta_m) \approx \max(L_m)$. Typically, $n_0 = 20$. In other words, the distance of the strong interactions must satisfy

$$R_{\text{Strong}} > \frac{n_0 \max(L_m)^2}{2\lambda}.$$
 (A6)

Since an edge shares two facets, a pair of edges (m, n) implies four facets. As shown in Fig. 13, assuming that the meshed triangles are identical, the sum over the facets of $Z_{m,n}$ can be made analytically. Since, for a given edge, $\delta_{m,n}^- = -\delta_{m,n}^+$ and $\rho_{m,n}^{p-,q-} = -\rho_{m,n}^{p+,q+}$, the sum over the four triangles of $s_p s_q e^{j(\phi_1+\phi_2)}$ reduces to

$$e^{j(\phi_1+\phi_2)} - e^{j(-\phi_1+\phi_2)} - e^{j(\phi_1-\phi_2)} + e^{-j(\phi_1+\phi_2)}$$

= $-4\sin\phi_1\sin\phi_2$. (A7)

In the same way, the sum over the four triangles of $s_p s_q e^{j(\phi_1 + \phi_2)} \rho_n^p \cdot \rho_n^q / (\rho_m^{p+} \cdot \rho_n^{q+})$ reduces to

$$e^{j(\phi_1+\phi_2)} + e^{j(-\phi_1+\phi_2)} + e^{j(\phi_1-\phi_2)} + e^{-j(\phi_1+\phi_2)}$$

= 4 \cos \phi_1 \cos \phi_2. (A8)

An element $\bar{Z}_{m,n}$ of the impedance matrix is then

$$\tilde{Z}_{m,n} \approx \frac{e^{-jkR_{m,n}}}{\pi R_{m,n}L_mL_n} \Big(\boldsymbol{\rho}_m^{p+} \cdot \boldsymbol{\rho}_n^{q+} \cos \phi_1 \cos \phi_2 + \frac{\sin \phi_1 \sin \phi_2}{k^2} \Big).$$
(A9)

Applying the following identities [23]:

$$\begin{cases} \cos(z\cos\theta) = J_0(z) + 2\sum_{p=1}^{\infty} J_{2p}(z)\cos(2p\theta) \\ \sin(z\cos\theta) = 2\sum_{p=1}^{\infty} (-1)^p J_{2p+1}(z)\cos[(2p+1)\theta] \end{cases}$$
(A10)

where J_p is the Bessel function of the first kind and order p, we have

$$\begin{cases} \cos \phi_1 \cos \phi_2 \approx J_0(q_m) J_0(q_n) \\ \sin \phi_1 \sin \phi_2 \approx -4 J_1(q_m) J_1(q_n) \cos \phi_m \cos \phi_n \end{cases}$$
(A11)

where $\phi_1 = -q_m \cos \phi_m$ and $\phi_2 = q_n \cos \phi_n$. The use of (A5) leads to

$$q_{m,n} = k\delta_{m,n} \left(1 - \frac{j}{kR_{m,n}} \right), \quad \cos\phi_{m,n} = \hat{\boldsymbol{\delta}}_{m,n} \cdot \hat{\boldsymbol{R}}_{m,n}.$$
(A12)

In (A10), only the first term of the sum is kept since $|q_{m,n}| \ll 1$.

In (A12), it is important to keep in mind that the extra term $j/(kR_{m,n})$ comes from a Taylor series expansion up to the order one. Then, we can simplify (A11) as

$$\cos\phi_{1}\cos\phi_{2} \approx J_{0}(w_{m})J_{0}(w_{n}) + \frac{j}{kR_{m,n}}$$

$$\times [w_{m}J_{1}(w_{m})J_{0}(w_{n}) + w_{n}J_{1}(w_{n})J_{0}(w_{m})]$$

$$\approx \left(1 - \frac{w_{m}^{2}}{4}\right)\left(1 - \frac{w_{n}^{2}}{4}\right) + \frac{j}{2kR_{m,n}}\left(w_{m}^{2} + w_{n}^{2}\right) \quad (A13)$$

and

$$\frac{\sin \phi_1 \sin \phi_2}{\cos \phi_m \cos \phi_n} \approx -4J_1(w_m)J_1(w_n) + \frac{2j}{kR_{m,n}}$$

$$\times \left\{ w_m J_1(w_n)[J_0(w_m) - J_2(w_m)] + w_n J_1(w_m)[J_0(w_n) - J_2(w_n)] \right\}$$

$$\approx -4J_1(w_m)J_1(w_n) + \frac{2j}{kR_{m,n}}$$

$$\times [w_m J_1(w_n)J_0(w_m) + w_n J_1(w_m)J_0(w_n)]$$

$$\approx -w_m w_n \left(1 - \frac{2j}{kR_{m,n}} \right)$$
(A14)

where $w_{m,n} = k \delta_{m,n}$. In addition, since $|w_{m,n}| \ll 1$, where $J_0(w_{m,n}) \approx 1 - w_{m,n}^2/4$ and $J_1(w_{m,n}) \approx w_{m,n}/2$. These approximations make the programming easier.

In conclusion, an element of the impedance can be approximated as

$$L_{m}L_{n}\tilde{Z}_{m,n} \approx G_{m,n}^{(1)} \left(1 - \frac{w_{m}^{2}}{4}\right) \left(1 - \frac{w_{n}^{2}}{4}\right) \rho_{m}^{p+} \cdot \rho_{n}^{q+} + \left(w_{m}^{2} + w_{n}^{2}\right) G_{m,n}^{(2)} \rho_{m}^{p+} \cdot \rho_{n}^{q+} + w_{m}w_{n}G_{m,n}^{(3)} \left(\boldsymbol{R}_{m,n} \cdot \hat{\boldsymbol{\delta}}_{m}\right) \left(\boldsymbol{R}_{m,n} \cdot \hat{\boldsymbol{\delta}}_{n}\right)$$
(A15)

where

$$\begin{cases} w_{m,n} = k \delta_{m,n} = k \| \boldsymbol{\delta}_{m,n} \| \\ \boldsymbol{\delta}_{m,n} = \boldsymbol{P}_{m,n} - \boldsymbol{G}_{p+,q+} \\ \boldsymbol{R}_{m,n} = \boldsymbol{P}_{m} - \boldsymbol{P}_{n} \\ \boldsymbol{\rho}_{m,n}^{p+,q+} = (\boldsymbol{V}_{m,n}^{p+,q+} - \boldsymbol{G}_{p+,q+})/2 \end{cases}$$
(A16)

and

$$\begin{cases}
G_{m,n}^{(1)} = \frac{e^{-jkR_{m,n}}}{\pi R_{m,n}} \\
G_{m,n}^{(2)} = G_{m,n}^{(1)} \frac{j}{2kR_{m,n}} \\
G_{m,n}^{(3)} = -G_{m,n}^{(1)} \left(1 - \frac{2j}{kR_{m,n}}\right) \frac{1}{k^2 R_{m,n}^2}
\end{cases}$$
(A17)

Since, for a planar surface, $\mathbf{R}_{m,n} = \mathbf{r}_{m,n} = (x_n - x_m)\hat{\mathbf{x}} + (y_n - y_m)\hat{\mathbf{y}}$ on a uniform grid, the matrix $\bar{\mathbf{Z}}_{m,n}^{(i)}$ associated with the element $G_{m,n}^{(i)}$ is Toeplitz. Then, the matrix product $O_m \bar{\mathbf{Z}}_{m,n}^{(i)} S_n$ can be computed from FFTs, where O_m and S_n are any observation and source vectors, respectively. For instance, the first term of (A15) is expanded as

$$(f_m \rho_{m,x}^{p+}) G_{m,n}^{(1)} (\rho_{n,x}^{q+} f_n) + (f_m \rho_{m,y}^{p+}) G_{m,n}^{(1)} (\rho_{n,y}^{q+} f_n) + (f_m \rho_{m,z}^{p+}) G_{m,n}^{(1)} (\rho_{n,z}^{q+} f_n)$$
(A18)

where $f_{m,n} = 1 - w_{m,n}^2/4$ and the subscripts (x, y, z) stand for the components of the vector. The above equation is in a form such that the source point is on the right-hand side of Green's function, while the observation point is located on the left-hand side of Green's function. In addition, three matrix– vector products are required, which are computed from FFTs. The same way is used for the last two terms of (A15), which requires six and 36 matrix–vector products, respectively, which are computed from FFTs. Thus, 45 matrix–vector products are necessary.

If the surface is not planar, then $R_{m,n} = \sqrt{r_{m,n}^2 + (z_n - z_m)^2}$, where z(x, y) is the elevation of the surface. A Taylor series expansion over $z = z_n - z_m$ up to the fourth order leads to

$$G_{m,n}^{(p)} = G_{m,n}^{(p),(0)} \left(1 + z^2 A_{m,n}^{(p),(1)} + z^4 A_{m,n}^{(p),(2)} \right)$$
(A19)

$$\begin{cases} A_{m,n}^{(1),(1)} = -\frac{1+ju}{2r_{m,n}^2} \\ A_{m,n}^{(2),(1)} = -\frac{2+ju}{2r_{m,n}^2} \\ A_{m,n}^{(3),(1)} = -j\frac{u^2-5ju-8}{2r_{m,n}^2(u-2j)} \end{cases}$$
(A20)
$$\begin{cases} A_{m,n}^{(3),(1)} = \frac{3+3ju-u^2}{8r_{m,n}^4} \\ A_{m,n}^{(2),(2)} = \frac{8+5ju-u^2}{8r_{m,n}^4} \\ A_{m,n}^{(2),(2)} = \frac{9ju^2+33u-48j-u^3}{8r_{m,n}^4(u-2j)} \end{cases}$$
(A21)

 $G_{m,n}^{(p),(0)} = G_{m,n}^{(p)}|_{z_n-z_m=0} = G_{m,n}^{(p)}(r_{m,n}), \text{ and } u = k \| \mathbf{r}_{m,n} \| = kr_{m,n} (\mathbf{r}_{m,n} = (x_n - x_m)\hat{\mathbf{x}} + (y_n - y_m)\hat{\mathbf{y}}). \text{ Since } \{A_{m,n}^{(p),(1)}, A_{m,n}^{(p),(2)}\} \text{ depends only on } r_{m,n}, \text{ on a uniform grid, the elements } \{G_{m,n}^{(p),(0)}A^{(p),(1)}, G_{m,n}^{(p),(0)}A^{(p),(2)}\} \text{ are also Toeplitz.}$

For the first order, since $R_{m,n} = r_{m,n}[1 + z^2/(2r_{m,n}^2)]$, the second order is neglected in comparison to the order zero if $kz^2/(2r_{m,n}) < 2\pi/n_0$ (condition on the phase of $e^{-jkR_{m,n}}$) and if $z^2/(2r_{m,n}^2) < 1/n_0$ (condition on the amplitude of $1/R_{m,n}$), where $n_0 > 1$ is an integer ranging from 10 to 20. Typically, $n_0 = 20$. This is consistent with (A20) and (A21). In other words, the strong interaction distance must satisfy

$$r_{\text{Strong}} > \max\left(\frac{\max(\Delta z)^2 n_0}{2\lambda}, \max(\Delta z)\sqrt{\frac{n_0}{2}}\right)$$
 (A22)

where $\Delta z = \max(z) - \min(z)$.

For a nonplanar surface, the conditions $\delta_{m,n}^{-} = -\delta_{m,n}^{+}$ and $\rho_{m,n}^{p^{-,q^{-}}} = -\rho_{m,n}^{p^{+,q^{+}}}$ are not satisfied. Then, the mean value is taken, that is, $\delta_{m,n}^{+} \rightarrow [\delta_{m,n}^{+} + (-\delta_{m,n}^{-})]/2$ and $\rho_{m,n}^{p^{+,q^{+}}} \rightarrow [\rho_{m,n}^{p^{+,q^{+}}} + (-\rho_{m,n}^{p^{-,q^{-}}})]/2$.

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