A Three-Dimensional Locally One-Dimensional Multiresolution Time-Domain Method Using Daubechies Scaling Function

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Abstract

A three-dimensional locally one-dimensional multiresolution time-domain(LOD-MRTD) method is introduced and unconditional stability is proved analytically. The updating formulations have fewer terms on the right-hand side than those of an alternating direction implicit MRTD(ADI-MRTD). The validation of the method is presented using the resonance frequency problem of an empty cavity. The reduction of the numerical dispersion technique is also combined with the proposed method. The numerical examples show that the combined method can improve the accuracy significantly.

Key words : Locally One-Dimensional Multiresolution Time-Domain(LOD-MRTD), Alternating Direction Implicit Multiresolution Time-Domain(ADI-MRTD), Numerical Dispersion, Artificial Anisotropy.

I. Introduction

The multiresolution time-domain(MRTD) method applying wavelets expansions in space has been developed for solving electromagnetic problems^[1]. Its numerical dispersion errors are significantly improved in comparison with the finite-difference time-domain(FDTD) method^[2]. However, the Courant-Friedich-Levy(CFL) stability condition still remains. An alternating direction implicit MRTD(ADI-MRTD)^{[3],[4]} method was proposed in order to overcome the CFL restriction. Recently, the unconditionally stable locally one-dimensional(LOD) time stepping technique has been applied to the FDTD method ^{[5]~[8]}. LOD-FDTD requires fewer arithmetic operations than ADI-FDTD while providing comparable accuracy.

As in the ADI method, the LOD method exhibits a splitting error that depends on not only the time step size, but also the spatial resolution. The time step is free of the CFL stability condition, but it cannot be made too large with the coarse mesh if a reasonable computational accuracy is to be achieved. To improve the accuracy, modified methods such as iterative techniques^{[9],[10]} and artificial anisotropy techniques^{[11]~[14]} were developed. The artificial anisotropy method takes only slightly more pre-computing than the conventional method, while the iterative method is computationally more expensive than

the conventional method. In this paper, the LOD technique is applied to the MRTD, resulting in an unconditionally stable 3-D LOD-MRTD method. The algorithm based on Daubechies scaling functions is only considered for simplicity. Artificial anisotropy is also extended to the 3-D LOD-MRTD method to improve the accuracy. Enhanced numerical accuracy is shown in several numerical results.

II. Formulations of 3-D LOD-MRTD Method

The LOD-MRTD can be derived by applying the LOD algorithm to the MRTD schemes. The procedure is similar to the LOD-FDTD developed in [8]. The shifted Daubechies scaling functions with two vanishing moments(D_2) are considered for a simple representation of the LOD-MRTD formulation. Maxwell's curl equations in isotropic and lossless media are given as

$$\nabla \times H = \varepsilon \frac{\partial E}{\partial t}$$
 $\nabla \times E = -\mu \frac{\partial H}{\partial t}$. (1)

These equations can be rewritten in the Cartesian coordinates as

$$\frac{\partial F}{\partial t} = [A]F + [B]F, \qquad (2)$$

where $F = \begin{bmatrix} E_x, E_y, E_z, H_x, H_y, H_z \end{bmatrix}^T$ and

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$$[A] = \begin{bmatrix} 0 & 0 & 0 & 0 & 0 & \frac{\partial}{\partial \partial y} \\ 0 & 0 & 0 & \frac{\partial}{\partial \partial z} & 0 & 0 \\ 0 & 0 & 0 & 0 & \frac{\partial}{\partial \partial x} & 0 \\ 0 & \frac{\partial}{\mu \partial z} & 0 & 0 & 0 & 0 \\ 0 & 0 & \frac{\partial}{\mu \partial x} & 0 & 0 & 0 \\ \frac{\partial}{\mu \partial y} & 0 & 0 & 0 & 0 & 0 \end{bmatrix},$$
$$[B] = \begin{bmatrix} 0 & 0 & 0 & 0 & \frac{\partial}{\partial z} & 0 \\ 0 & 0 & 0 & 0 & \frac{\partial}{\partial z} & 0 \\ 0 & 0 & 0 & \frac{\partial}{\partial z} & 0 \\ 0 & 0 & 0 & \frac{\partial}{\partial z} & 0 & 0 \\ 0 & 0 & \frac{\partial}{\mu \partial y} & 0 & 0 & 0 \\ \frac{\partial}{\mu \partial z} & 0 & 0 & 0 & 0 \\ 0 & \frac{\partial}{\mu \partial x} & 0 & 0 & 0 & 0 \end{bmatrix}$$

in which ε and μ present permittivity and permeability, respectively. The Crank-Nicolson(CN) approximation is applied and then (2) can be split into the following two sub-steps

$$\left(I - \frac{\Delta t}{2} [A]\right) F^{n+\frac{1}{2}} = \left(I + \frac{\Delta t}{2} [A]\right) F^{n}$$

$$\left(I - \frac{\Delta t}{2} [B]\right) F^{n+1} = \left(I + \frac{\Delta t}{2} [B]\right) F^{n+\frac{1}{2}},$$
(3b)

where *I* denotes the unit matrix. The Yee's grid is employed to discretize the field space, and the field quantities are only expanded in terms of the scaling function expansions. Therefore, the two sub-step procedures can be formulated as follows:

First sub-step from n to n+1/2Implicit updating for E_x

$${}_{x}E_{i+1/2,j,k}^{n+1/2} - \frac{\left(\Delta t\right)^{2}}{4\mu\varepsilon\left(\Delta y\right)^{2}} \sum_{m=-3}^{2} a(m) \sum_{\nu=-3}^{2} a(\nu)_{x}E_{i+1/2,j+m+\nu+1,k}^{n+1/2}$$

$$= {}_{x}E_{i+1/2,j,k}^{n} + \frac{\left(\Delta t\right)^{2}}{4\mu\varepsilon\left(\Delta y\right)^{2}} \sum_{m=-3}^{2} a(m) \sum_{\nu=-3}^{2} a(\nu)_{x}E_{i+1/2,j+m+\nu+1,k}^{n}$$

$$+ \frac{\Delta t}{\varepsilon\Delta y} \sum_{m=-3}^{2} a(m)_{z}H_{i+1/2,j+m+1/2,k}^{n}$$
(4)

Explicit updating for H_z

$$z H_{i+1/2,j+1/2,k}^{n+1/2} = z H_{i+1/2,j+1/2,k}^{n+1/2} + \frac{\Delta t}{2\mu\Delta y} \sum_{\nu=-3}^{2} a(\nu) \Big({}_{x} E_{i+1/2,j+\nu+1,k}^{n+1/2} - {}_{x} E_{i+1/2,j+\nu+1,k}^{n} \Big)$$
(5)

Second sub-step from n+1/2 to n+1Implicit updating for E_x

$${}_{x}E_{i+1/2,j,k}^{n+1} - \frac{\left(\Delta t\right)^{2}}{4\mu\varepsilon\left(\Delta z\right)^{2}} \sum_{m=-3}^{2} a(m) \sum_{\nu=-3}^{2} a(\nu)_{x}E_{i+1/2,j,k+m+\nu+1}^{n+1}$$

$$= {}_{x}E_{i+1/2,j,k}^{n+1/2} + \frac{\left(\Delta t\right)^{2}}{4\mu\varepsilon\left(\Delta z\right)^{2}} \sum_{m=-3}^{2} a(m) \sum_{\nu=-3}^{2} a(\nu)_{x}E_{i+1/2,j,k+m+\nu+1}^{n+1/2}$$

$$- \frac{\Delta t}{\varepsilon\Delta z} \sum_{m=-3}^{2} a(m)_{y}H_{i+1/2,j,k+m+1/2}^{n+1/2}$$
(6)

Explicit updating for H_v

$${}_{v}H_{i+1/2,j,k+1/2}^{n+1} = {}_{v}H_{i+1/2,j,k+1/2}^{n+1/2} - \frac{\Delta t}{2\mu\Delta z} \sum_{\nu=-3}^{2} a(\nu) \Big({}_{x}E_{i+1/2,j,k+\nu+1}^{n+1} - {}_{x}E_{i+1/2,j,k+\nu+1}^{n+1/2} \Big),$$
(7)

where the coefficients a(v) (or a(m)) are the connection coefficients of the scaling function expansions. The coefficients a(v) are reported in [4] and shown in Table 1. The values of a(v) for v<0 are given by a(-1-v) = -a(v). Δx , Δy and Δz represent the space discretization intervals in x-, y- and z-direction, and Δt is the time interval. The complete formulations for the unconditionally stable LOD-MRTD scheme can be obtained by applying the same procedure for other components. For derived formulations, there are six implicit and six explicit equations to be computed in one time step, which is the same number as those of the conventional ADI-MRTD method^[3]. However, as shown in (4) \sim (7), the right-hand sides(RHS) of the equations have fewer terms than those in the ADI-MRTD method. The field component E_x , the second term of the RHS in (4), can be calculated at eleven adjacent grid points, while the field component is computed at 36 neighbor grid points in ADI-MRTD. Further, the introduced scheme needs only the summation of the field component H_z , the third term of RHS in (4), while the ADI-MRTD method needs the summations of both H_z and H_y . Clearly, the LOD-MRTD computation is more concise than the ADI-MRTD.

III. Numerical Stability

The equations in the spatial spectral domain can be written in the following matrix forms.

Table	1.	Coefficient	a((v))
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v	<i>a</i> (<i>v</i>)
0	1.22916661202745
1	-0.09374997764764
2	0.01041666418309

Sub-step 1:
$$F^{n+1/2} = \Lambda_1 F^n$$
 (8a)

Sub-step 2:
$$F^{n+1} = \Lambda_2 F^{n+1/2}$$
, (8b)

where

$$\Lambda_{1} = \begin{bmatrix} \frac{G_{y}}{Q_{y}} & 0 & 0 & 0 & 0 & \frac{-j2W_{y}}{\varepsilon Q_{y}} \\ 0 & \frac{G_{z}}{Q_{z}} & 0 & \frac{-j2W_{z}}{\varepsilon Q_{z}} & 0 & 0 \\ 0 & 0 & \frac{G_{x}}{Q_{x}} & 0 & \frac{-j2W_{x}}{\varepsilon Q_{x}} & 0 \\ 0 & 0 & \frac{-j2W_{z}}{\mu Q_{z}} & 0 & \frac{G_{z}}{Q_{z}} & 0 & 0 \\ 0 & 0 & \frac{-j2W_{x}}{\mu Q_{x}} & 0 & \frac{G_{x}}{Q_{x}} & 0 \\ \frac{-j2W_{y}}{\mu Q_{y}} & 0 & 0 & 0 & 0 & \frac{G_{y}}{Q_{y}} \end{bmatrix}$$

$$Q_{\alpha} = 1 + \frac{W_{\alpha}^{2}}{\mu \varepsilon}, \quad G_{\alpha} = 1 - \frac{W_{\alpha}^{2}}{\mu \varepsilon}, \quad \alpha = x, y, z$$

$$W_{\alpha} = \frac{\Delta t}{\Delta \alpha} \sum_{m=0}^{2} a(m) \sin \left(\frac{k_{\alpha}^{\ell} \Delta \alpha}{2} (2m+1) \right)$$

$$\Lambda_{2} = \begin{bmatrix} \frac{G_{z}}{Q_{z}} & 0 & 0 & 0 & \frac{j2W_{z}}{\varepsilon Q_{x}} & 0 \\ 0 & 0 & \frac{G_{y}}{Q_{y}} & \frac{j2W_{y}}{\varepsilon Q_{y}} & 0 & 0 \\ 0 & 0 & \frac{j2W_{y}}{\mu Q_{y}} & \frac{G_{y}}{Q_{y}} & 0 & 0 \\ \frac{j2W_{z}}{\mu Q_{z}} & 0 & 0 & 0 & \frac{G_{z}}{Q_{z}} & 0 \\ 0 & \frac{j2W_{x}}{\mu Q_{x}} & 0 & 0 & 0 & \frac{G_{x}}{Q_{x}} \end{bmatrix}$$

The above two step equations can be written as

$$F^{n+1} = \Lambda_2 \cdot \Lambda_1 F^n = \Lambda F^n \,. \tag{9}$$

The eigenvalues of the Λ can be found as

$$\begin{split} \lambda_1 &= \lambda_2 = 1 \\ \lambda_3 &= \lambda_4 = x + jy \\ \lambda_5 &= \lambda_6 = x - jy \end{split}$$

where

$$x = \frac{2(Q_x + Q_y + Q_z - Q_xQ_y - Q_yQ_z - Q_zQ_x) + Q_xQ_yQ_z}{Q_xQ_yQ_z}$$
$$y = \sqrt{1 - x^2} .$$

It is easy to show that $|\lambda_3| = |\lambda_4| = |\lambda_5| = |\lambda_6| = 1$. Therefore, the introduced LOD-MRTD method is unconditionally stable.

IV. Numerical Examples

For validation of the proposed LOD-MRTD method, an empty cavity with dimensions of 90×60×150 mm was computed. In the first case, the mesh chosen was rather dense. For a uniform standard Yee's cell, the spatial steps in all dimensions were chosen to be $\Delta \alpha = 10$ mm with $\alpha = x$, y and z, leading to a 9×6×15 grid. The time step was varied from 0.2 to five times the CFL limit of the original scaling function based MRTD(S-MRTD). The CFL time step limit of the original S-MRTD with D₂ is $0.433(\Delta \alpha/c_0)$, where c_0 is the speed of light in the vacuum. Electric fields are sampled for a period of 108.3 ns. Namely, the iterations are assumed to be 5,000 when the relative time step is 0.5. Fig. 1 shows the relative errors from the analytical solutions for the two resonant modes, TE₁₀₁ and TE₁₀₂. The horizontal axis is the ratio of the LOD-MRTD time step Δt to the CFL time step limit ($\Delta t_{MRTD-CFL}=0.433(\Delta \alpha/c_0)$). It can be seen that, when the time step increases to four times the CFL time step limit, the errors are still relatively small, less than 2.78 % for the TE_{101} mode. In the second case, a coarse mesh was chosen with the space step being 30 mm. It led to a 3×2×5 grid. Fig. 2 shows the computation results. With a coarse mesh, the time step can no longer be too large. Even with the time step of twice the CFL time step limit, the error with the LOD-MRTD for the lowest resonant mode (TE_{101}) becomes lager than 5.61 %. When the relative time step is 0.5, Table 2 shows the numerical results in terms of the mesh size, lowest resonance frequency, relative



Fig. 1. Relative resonance frequency errors in the cavity $(9 \times 6 \times 15 \text{ cells})$ as a function of the relative time step(TE₁₀₁: 1.942 GHz, TE₁₀₂: 2.602 GHz).



Fig. 2. Relative resonance frequency errors in the cavity $(3 \times 2 \times 5 \text{ cells})$ as a function of the relative time step(TE₁₀₁: 1.942 GHz, TE₁₀₂: 2.602 GHz).



Fig. 3. Comparison of the relative resonance frequency errors in the cavity($3 \times 2 \times 5$ cells) for the lowest resonance frequency(TE₁₀₁: 1.942 GHz).

error, and computational time. Frequency is given in GHz, error in %, and time in seconds. These simulations are performed on Intel core 2 duo (2.33 GHz) PC. As can be seen, the errors for both ADI- and LOD-MRTD are at the same level, but the computational time is less in LOD-MRTD. The FDTD method needs denser spatial intervals in order to obtain the accurate results. The relative errors of the LOD- MRTD and the ADI-MRTD from the analytical solutions, with respect to relative time steps, are shown in Fig. 3. It is found that both methods have the same relative errors. Although the time step is free of the CFL stability condition, it cannot be made too large in case of the coarse mesh.

Table	2	Comparison	of	numerical	results
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	FDTD		ADI-MRTD		LOD-MRTD	
Mesh (Iteration)	Results (GHz)	Time	Results (GHz)	Time (sec)	Results (GHz)	Time (sec)
	(Error)	(sec)	(Error)		(Error)	
3×2×5 (5,000)	1.875	2.25	1.941	46.76	1.941	16.98
	(3.40 %)	5.25	(0.05 %)		(0.05 %)	
6×4×10 (10,000)	1.925	13.82				
	(0.82 %)					
9×6×15 (15,000)	1.935	36.70				
	(0.31 %)					

V. Numerical Dispersion and Optimization

The numerical dispersion relation can be established using a procedure similar to the one described in [15]. Equation (9) is rewritten in the spectral domain as

$$\left(e^{j\omega\Delta t}I - \Lambda\right)F = 0 \tag{10}$$

The resulting numerical dispersion relation for freespace is

$$\tan^{2}\left(\frac{\omega\Delta t}{2}\right) = \frac{P_{x} + P_{y} + P_{z} + P_{x}P_{y} + P_{y}P_{z} + P_{z}P_{x}}{1 + P_{x}P_{y}P_{z}},$$
 (11)

where

$$P_{\alpha} = \frac{W_{\alpha}^2}{\mu_0 \varepsilon_0}, \qquad \alpha = x, y, z$$

In order to reduce the dispersion error, the numerical dispersion relation in magnetically and electrically anisotropic material can be considered. Assume that the relative permeability is equal to the relative permittivity (i.e., ε_r =diag(ε_x , ε_y , ε_z)= μ_r). The numerical dispersion relation can be obtained as the following form in free space.

$$\tan^{2}\left(\frac{\omega\Delta t}{2}\right) = \left[\frac{P_{x}}{\varepsilon_{y}\varepsilon_{z}} + \frac{P_{y}}{\varepsilon_{z}\varepsilon_{x}} + \frac{P_{z}}{\varepsilon_{x}\varepsilon_{y}} + \frac{P_{x}P_{y}}{\varepsilon_{x}\varepsilon_{y}\varepsilon_{z}^{2}} + \frac{P_{y}P_{z}}{\varepsilon_{x}^{2}\varepsilon_{y}\varepsilon_{z}} + \frac{P_{z}P_{x}}{\varepsilon_{x}\varepsilon_{y}^{2}\varepsilon_{z}}\right] / \left[1 + \frac{P_{x}P_{y}P_{z}}{\varepsilon_{x}^{2}\varepsilon_{y}^{2}\varepsilon_{z}^{2}}\right]$$
(12)

It is exactly the same as the relation of the ADI-MRTD method. Therefore, the reduction procedure for the numerical dispersion error can be employed in the LOD-MRTD method, as it previously worked for the ADI-MRTD method^[14]. As in ADI-MRTD, a Courant coefficient *S* and the spatial resolution coefficient *R* are defined as

$$S = c_0 \Delta t \sqrt{\frac{1}{\varepsilon_y \varepsilon_z \Delta x^2} + \frac{1}{\varepsilon_z \varepsilon_x \Delta y^2} + \frac{1}{\varepsilon_x \varepsilon_y \Delta z^2}}$$
(13)

$$R = \frac{\lambda}{\sqrt{\Delta x^2 + \Delta y^2 + \Delta z^2}} \,. \tag{14}$$

The parameters are introduced to denote the shape of a single Yee's cell.

$$Z_1 = \frac{\Delta x}{\Delta y}, \quad Z_2 = \frac{\Delta x}{\Delta z}, \quad Z = \sqrt{1 + Z_1^{-2} + Z_2^{-2}},$$
 (15)

where $\lambda = c_0/f$ is the wavelength and c_0 is the speed of light in the vacuum. If it is supposed that the wave propagation at angles ϕ and Θ is in the spherical coordinate system, then $\mathcal{K}_x = 7 \sin\Theta \cos \phi$, $\mathcal{K}_y = 7 \sin\Theta \sin \phi$, $\mathcal{K}_z =$ $\mathcal{K} \cos\Theta$, and \mathcal{K} is denoted as $2\pi/\tilde{\lambda}$, where $\tilde{\lambda}$ is the wavelength in the grid. A coefficient $A = (v_p/c_0) = (\tilde{\lambda}/\lambda)$ is defined as the normalized numerical phase velocity, where v_p is the propagation velocity in the grid. In order to find the optimization parameter values for ε_x , ε_y , and ε_z that ensure the minimum numerical dispersion, the values are first determined to yield A=1 along the coordinate directions ($\phi=0^\circ$, $\Theta=90^\circ$), ($\phi=90^\circ$, $\Theta=90^\circ$), and ($\Theta=0^\circ$). The optimization parameter values for ε_x , ε_y , and ε_z are calculated as follows:

$$\varepsilon_x = \frac{RZZ_1Z_2L}{\pi S} \sqrt{\left(\frac{U_2U_3}{U_1}\right)} \cdot \tan^{-1}\left(\frac{S}{L}\right)$$
(16)

$$\varepsilon_{y} = \frac{RZZ_{2}L}{\pi SZ_{1}} \sqrt{\left(\frac{U_{1}U_{3}}{U_{2}}\right)} \cdot \tan^{-1}\left(\frac{S}{L}\right)$$
(17)

$$\varepsilon_{z} = \frac{RZZ_{1}L}{\pi SZ_{2}} \sqrt{\left(\frac{U_{1}U_{2}}{U_{3}}\right)} \cdot \tan^{-1}\left(\frac{S}{L}\right), \qquad (18)$$

where

$$U_{1} = \left\{ \sum_{m=0}^{2} a(m) \sin\left[\frac{\pi (2m+1)}{ARZ}\right] \right\}^{2}$$
$$U_{2} = \left\{ \sum_{m=0}^{2} a(m) \sin\left[\frac{\pi (2m+1)}{ARZZ_{1}}\right] \right\}^{2}$$
$$U_{3} = \left\{ \sum_{m=0}^{2} a(m) \sin\left[\frac{\pi (2m+1)}{ARZZ_{2}}\right] \right\}^{2}$$
$$L = \sqrt{U_{1}^{-1} + U_{2}^{-1} + U_{3}^{-1}}$$

After the calculated optimization parameters are applied to update the schemes, the relative numerical phase velocity A is 1 along the axial directions. The phase velocity anisotropic error is obtained as

$$\Delta \mathscr{H}_{aniso} = \frac{\max\left(v_p(\phi, \theta)\right) - \min\left(v_p(\phi, \theta)\right)}{c_0}$$



Fig. 4. Normalized phase velocity versus wave angles for standard and optimized LOD-MRTD.



Fig. 5. Relative resonance frequency errors in the cavity (90×60×150 mm) for standard and optimized LOD-MRTD(TE₁₀₁: 1.942 GHz).

When min $(v_p(\phi, \Theta))$ is equal to 1, it is possible to obtain the optimization parameters by setting $A=1-\Delta \tilde{v}$ aniso/2. On the other hand, setting $A=1+\Delta \tilde{v}_{aniso}/2$ yields the optimization parameters, while max $(v_p(\phi, \Theta))$ is equal to 1. Under this optimization, the numerical phase velocity errors have their minimum values in all propagation directions. Fig. 4 shows the effect of the correction with wave propagation angle. The horizontal top axis is the wave propagation angle Θ , while the angle ϕ is 45°. When the angle θ is 90°, the horizontal bottom axis is the wave propagation angle ϕ . A spatial resolution coefficient R=2.9709 and the Courant coefficient $S=1.125(\Delta t \approx 1.45 \Delta t_{MRTD-CFL})$ are considered. The parameters are optimized at 1.942 GHz. The calculated correction parameters are $\varepsilon_x = \varepsilon_y = \varepsilon_z = 0.9667$ for the uniform grid. The maximum standard dispersion error is -4.16 %, whereas the corrected dispersion error

	The optimized LOD-MRTD scheme						
Analytical result (GHz)	$\Delta t=0.987$ $\Delta t_{MRTD-CFL}$		$\Delta t=1.88$ $\Delta t_{MRTD-CFL}$		$\Delta t=2.62$ $\Delta t_{\rm MRTD-CFL}$		
	Results (GHz)	Time	Results (GHz)	Time	Results (GHz)	Time	
	(Error)	(sec)	(Error)	(sec)	(Error)	(sec)	
1.942	1.943	8.61	1.950	4.55	1.962	3.30	
	(0.05 %)		(0.41 %)		(1.03 %)		

Table 3. Numerical results with different Δt .

 Table 4. Comparison of results with standard and optimized LOD-MRTD.

$\Delta t / \Delta t_{MRTD}$ -CFL	Standard LOI	D-MRTD	Optimized LOD-MRTD		
	Simulation results(MHz)	Relative error	Simulation results(MHz)	Relative error	
0.9854	123.27	1.38 %	125.05	0.04 %	
1.4433	121.01	3.19 %	125.43	0.34 %	
1.8644	118.36	5.31 %	125.95	0.76 %	
2.2438	115.61	7.51 %	126.59	1.27 %	

remains within ± 1 %. The cavity with dimensions 90×60 ×150 mm is also computed with the optimization. The uniform spatial step $\Delta \alpha$ is chosen to be 30 mm, which leads to $3 \times 2 \times 5$ cells. The correction parameters are optimized to wavelength $\lambda = 0.1544$ m(R=2.9709) at each time step. Fig. 5 illustrates the relative resonance frequency errors for the standard and corrected methods. When the relative time step is 2.62, the error becomes 1.03 % after the correction procedure, while the error without correction becomes larger than 9.2 %. Table 3 shows the lowest resonance frequency, relative error, and computational time in the optimized scheme. Their physical time periods are the same, but the number of iterations is different. When the time steps $\Delta t=0.987\Delta t$ MRTD-CFL=4.277×10⁻¹¹ s, Δt =1.88 $\Delta t_{MRTD-CFL}$ =8.147×10⁻¹¹ s, and $\Delta t = 2.62 \Delta t_{\text{MRTD-CFL}} = 1.137 \times 10^{-10}$ s are chosen, the iterations are 2532, 1329, and 952, respectively. Table 4 shows the lowest relative resonance frequency errors in the cavity with $1.5 \times 1 \times 2$ m. The uniform spatial step $\Delta \alpha$ is chosen to be 0.5 m, which leads to $3 \times 2 \times 4$ cells. The correction parameters are optimized to an analytical resonance frequency of 125 MHz(R=2.769) at each time step. It is obvious that using the combined method can improve the accuracy.

VI. Conclusion

The unconditionally stable 3-D MRTD scheme based on the LOD method is introduced. Stability restriction associated with the conventional S-MRTD method is completely removed. The numerical errors of this method are same as those of the ADI-MRTD method. However, it requires less computational expenditure than the ADI-MRTD method because the updating equations have fewer terms on their right-hand sides. Moreover, this paper shows that the introduced work can be combined with an artificial anisotropy technique in order to reduce the numerical dispersion error. The combined method shows more accurate numerical results, even though the time step increases to several times of CFL limit of S-MRTD. Therefore the introduced LOD- MR-TD method combined with dispersion reduction procedure can yield more efficient numerical computation for solving electromagnetic field problems.

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