Uncertainty Analysis of Reflection Coefficient for a Coating with Random Flaws Using Adaptive Mesh and DGTD Method

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Abstract — An imperfect coating shall cause uncertainties in the analysis of electromagnetic properties. To quantify the influence of irregularity, complexity, and uncertainty of the coatings for electronic devices, an adaptive mesh algorithm combined with the discontinuous Galerkin time domain (AM-DGTD) method is developed. The uncertain variations are incorporated into the proposed algorithm by an appropriate parameterization. The standard statistical analysis is performed to calculate the appropriate moments, i.e., mean and variance. The developed method is validated by modeling a dielectric coating with uncertain flaws in an adaptive mesh grid. The computed quantities of interest from numerical estimations are compared with the analytical values, these results agree with the physical explanation, and are in good agreement with the exact values, as demonstrated by numerical experiments.

Index Terms—AM-DGTD method, reflection coefficient, statistical analysis, uncertainty quantification.

I. INTRODUCTION

Dielectric coatings applied to surfaces of aircraft frames, automobile frames and ship hull are able not only to protect objects from sunlight, moisture, dust and abrasion but also to improve their appearance [1]. For example, cracks, voids, inclusions, structural flaws and material defects [2], these uncertain flaws may impact the coating’s electromagnetic performance and cause difference from the ideal manufacturing objectives, or to damage the essential functions of the dielectric coatings. In order to engage in the practical analysis and simulation, one must find a suitable, flexible and efficient method to deal with these electromagnetic problems.

In the recent years, the method of moments (MoM) with probabilistic technique outlined in [3] was applied to determine the uncertainty in practical EM compatibility measurements. Polynomial chaos technique was used in the finite difference time domain (FDTD) method to study microwave circuits and free-space scattering problems [4]-[5]. An efficient stochastic finite difference time domain (S-FDTD) method [6], which uses truncated Taylor series approximations in the derivation, is employed for evaluating statistical variation in the EM fields caused by variability or uncertainty in the electrical properties of the materials in the model. To obtain static response characteristics of graphite/epoxy composite laminates with random material properties, Navaneetha Raj et al. employed Monte Carlo simulation and the finite element method (FEM) with different boundary conditions [7]. A hybrid approach was adopted in the spectral stochastic finite element method (SSFEM) and polynomial chaos expansion (PCE) to provide response analysis of a linear structure with uncertainties in both the structural parameters and the external excitation [8]. In [9], a probabilistic approach based on high-order accurate expansions of general stochastic processes and high-order discontinuous Galerkin method was applied to solve the time-domain electromagnetic problems with uncertainty in initial conditions, boundary conditions, sources, materials, computational domain, and/or geometries. In [10], Li et al. proposed a hybrid method which integrated an adaptive hierarchical spare grid collocation (ASGC) method and the discontinuous Galerkin time-domain (DGTD) method. The ASGC method is employed to approximate the stochastic observables of interest using interpolation functions over a set of collocation points determined by the Smolyak’s algorithm integrated with an adaptive strategy. Owing to the random flaws in the coatings are generally smaller, irregular, and more arbitrary in nature. Moreover, the DGTD method offers high-order accuracy to a coarse resolution, geometrical flexibility through fully unstructured grids and higher computational efficiency, so it could be more suitable to be employed for the uncertainty quantization [9]-[14].

In this work, using the node displacement method and the node insertion method [15]-[17], we develop an
The node insertion method is simple, however, it is time and resource-consuming, because the insertion of DNV into \( V_x \) cause reduction in the minimum distance between adjacent nodes, then the smaller time step has to be employed to maintain the CFL condition. Pseudo algorithm for this adaptive mesh by inserting node is provided in Algorithm 1.

**Algorithm 1 Insertion algorithm for desired nodes**

**Input:** computational domain \( \Omega = [x_{min}, x_{max}] \), number of elements \( K \), desired node vectors \( DNV \)

**Output:** updated number of elements \( K^* \), updated global node vector \( V_x^* \)

```
for \( i \leftarrow 1 \) to \( K \) do
    /* Generate a simple equidistant grid and uniform node vectors \( V_x^* \) */
    \( V_x(i) \leftarrow (x_{max} - x_{min}) \cdot (i - 1) / K + x_{min} \)
    /* Insert the DNV into the uniform node vector \( V_x^* \) */
    /* Sort and return the unique values */
    \( V_x^* \leftarrow \text{unique}(\{V_x, DNV\}) \)
    \( K^* \leftarrow \text{length}(V_x^*) - 1 \)
Return \( K^*, V_x^* \)
```

**B. Node displacement mesh**

In this type of adaptive mesh generation, the adaptation can be achieved through suitably displacing the UNV. The algorithm identifies the nodes from the UNV which are nearest to the set of DNV and shifts these nodes to the desired locations. If a DNV coincides with one of the UNV, such as the node ① and ②, then no displacement is needed; if a structure represented by any two of the DNV (e.g., node ② and ③) is much larger than uniform elements, then the nearest uniform nodes are modified at the boundaries; if a structure is very small, such as the segment composed by node ④ and ⑤, the algorithm makes sure that at least one element is contained. Hence, only optimal numbers of element are generated, as shown in Fig. 2.

Note that the number of global nodes remains unchanged, that is, \( K^* = K \). This algorithm generates optimal density of mesh cells over the whole \( \Omega \).
regardless of geometry structure. The displacement of nodes to different position caused non-uniformity in the distribution of global nodes. Realization for this adaptive mesh algorithm is given below in Algorithm 2.

![Adaptive Nodes by Displacement Method](image)

Fig. 2. The mesh cell mappings for node displacing mesh algorithm.

**Algorithm 2 Displacement algorithm for desired nodes**

**Input**: computational domain \( \Omega = [x_{min}, x_{max}] \), number of elements \( K \), desired node vectors \( DNV \)

**Output**: updated number of elements \( K^* \), updated global node vector \( V_i^* \)

```plaintext
for i←1 to K do
    /* Generate a simple equidistant grid and uniform node vectors \( V_i \) */
    \( V_i(i) \leftarrow (x_{max} - x_{min}) \times (i-1)/K + x_{min} \)

if length(\( V \)) > 2*length(DNV)+2 then
    /*Generate index of DNV in set of the uniform node vectors UNV */
    [Val(q) Index(q)] ← min(abs(V_i-DNV(q)))
    /* Deal with nodes which near boundary of the domain */
    ConflictIndex ← find( Index(end)==Index)
    if length(ConflictIndex) > 1 then
        Index(end) ← Index(end)+1
    return

V_i (Index(q)) ← DNV(q)
V_i^* ← V_i
K^* ← length(V_i^*)-1
```

**Return**

\( K^*, V_i^* \)

C. Numerical scheme for one-dimensional case

To compute the reflection coefficients of the dielectric coating with random flaws, we solve the one-dimensional Maxwell’s equations in the time domain, subject to a broadband initial condition, and collect one time-trace at an observed point in the computational domain. Consider a lossless material TEM case, the time-dependent Maxwell’s equations can be written as follows:

\[
\varepsilon(z) \frac{\partial E_z}{\partial t} + \mu(z) \frac{\partial H_z}{\partial t} = 0, \quad \mu(z) \frac{\partial E_z}{\partial t} + \varepsilon(z) \frac{\partial H_z}{\partial t} = 0, \quad (1)
\]

where \( E_z \), \( H_z \), \( \varepsilon \) and \( \mu \) represent the electric field, the magnetic field, the local electric permittivity, and the local magnetic permeability, respectively.

Using the aforementioned adaptive mesh approaches, the computational domain \( \Omega = [l, r] \), which is tessellated by \( K^* \) subdomains, i.e., any one of elements of the physical space \( D^k \) is equal to \([z^k, z^k], k = 1...K^*\). The solution of Equation (1) will be discontinuous between elements. In an arbitrary element \( D^k \), the fields can be approximately expanded using local high-order Lagrange interpolation polynomial \( \ell^i_k(z) \):

\[
\left[ E^k_{z_1}(z,t) \right] = \left[ E^k_{z_1}(z,t) \right] = \sum_{n=1}^{N_p} \left[ E^k_{z_1}(z^k,t) \right] \ell^i_k(z), \quad (2)
\]

where \( N_p \) stands for the number of the local expansion. \( E^k_{z_1} \) and \( H^k_{z_1} \) contain a \( N_p \)-vector of expansion coefficients to be solved. \( \ell^i_k(z) \) signifies an \( N_p \)th order Lagrange polynomial. The relationship between \( N_p \) and the polynomial expansion order \( N \) is \( N_p = N + 1 \). On account of the fact that correctly choosing interpolation nodes can bring about good numerical behaviors, this work employs the Legendre-Gauss-Lobatto (LGL) interpolating nodes as \( z_i \) [21]-[23].

Next, multiplying (1) by a test function \( \ell^i_k(z) \) in an element \( D^k \), yields:

\[
\int_{D^k} \left[ \varepsilon(z) \frac{\partial E_z^k(z,t)}{\partial t} + \mu(z) \frac{\partial H_z^k(z,t)}{\partial t} \right] \ell^i_k(z) dz = 0,
\]

\[
\int_{D^k} \left[ \mu(z) \frac{\partial E_z^k(z,t)}{\partial t} + \varepsilon(z) \frac{\partial H_z^k(z,t)}{\partial t} \right] \ell^i_k(z) dz = 0. \quad (3)
\]

In order to couple with adjacent elements, Equation (3) are manipulated with integration by parts twice, and the strong variational formulation can be obtained as [21]:

\[
\int_{D^k} \left[ \varepsilon(z) \frac{\partial E_z^k}{\partial t} + \mu(z) \frac{\partial H_z^k}{\partial t} \right] \ell^i_k(z) dz = \int_{D^k} \hat{n} \left[ \frac{\partial H_z^k}{\partial z} - \frac{\partial E_z^k}{\partial z} \right] \ell^i_k(z) dz,
\]

\[
\int_{D^k} \left[ \mu(z) \frac{\partial H_z^k}{\partial t} + \varepsilon(z) \frac{\partial E_z^k}{\partial t} \right] \ell^i_k(z) dz = \int_{D^k} \hat{n} \left[ \frac{\partial E_z^k}{\partial z} - \frac{\partial E_z^k}{\partial z} \right] \ell^i_k(z) dz. \quad (4)
\]

Here, \( \hat{n} \) denotes the local outward pointing normal. On the right-hand side (RHS) of (4) \( (E_z^k, H_z^k) \) are numerical fluxes to exchange the coupling between neighboring elements. Using the Riemann conditions
and for stability reasons, we use a pure upwind flux [21] which could strongly damp unphysical modes,

\[
E^*_i = \frac{1}{|Y|} \left( \left\{ (YE^*_r) \right\} + \frac{1}{2} \left\{ [H^*_r] \right\} \right),
\]

\[
H^*_r = \frac{1}{|Z|} \left( \left\{ (ZH^*_i) \right\} + \frac{1}{2} \left\{ [E^*_i] \right\} \right),
\]

where \( Z = (Y)^{-1} = \sqrt{\mu / \varepsilon} \), represents the impedance of the medium.

Now substitute the expansions in (2) with the numerical flux of (5) into (4) and assume a smooth material in each element. After some algebraic computations, the explicit semi-discrete scheme in matrix-vector form can be obtained as follows:

\[
\frac{dE_{sh}^i}{dt} = \frac{1}{J^i c^i} \left\{-D^i H_{sh}^i + \frac{1}{M^i} \left[ \ell^i \left( z \right) \left( H_{sh}^i - H^*_r \right) \right] \right\},
\]

\[
\frac{dH_{sh}^i}{dt} = \frac{1}{J^i c^i} \left\{-D^i E_{sh}^i + \frac{1}{M^i} \left[ \ell^i \left( z \right) \left( E_{sh}^i - E^*_i \right) \right] \right\}.
\]

Here, the matrices \( J^i, D^i, \) and \( M^i \) represent the local transformation Jacobian, differentiation matrix and mass-integration matrix, respectively (see [21] for details).

The semi-discrete system of (6) is ordinary differential equations with respect to time. The forth-order low-storage explicit Runge-Kutta (LSERK) solver is employed for the time integration of (6) [24].

III. AM-DGTD FOR DETERMINISTIC PARAMETERS

A. Computation model

The objective of this article is to analyze the characteristics of a dielectric coating. To reduce the complexity of the problem, this coating is located at the center (i.e., \( x^{c0} = 0 \)) of the computational domain \( \Omega = [-6.0cm, 6.0cm] \). Free space is set on the both sides of the coating and the relative parameters \( \varepsilon = 1 \) and \( \mu = 1 \). The main sensitive variables with uncertainty are supposed to be mean values, i.e., the average thickness of the coating \( d = 2.0cm \) and its average relative permittivity \( \varepsilon = 4 \).

To model the excitation source, an \( x \)-polarized, \( z \)-directed Gaussian pulse with respect to space is used and the following initial conditions \((t=0)\) are adopted as:

\[
E_i = \exp \left(-4\pi \frac{(z-z_0)^2}{z_f^2}\right), \quad z_0 = -3.5cm, \quad z_f = 0.5cm,
\]

\[
H_i = 0
\]

where \( z_0 \) and \( z_f \) are the center position and the width of the Gauss curve, respectively. This Gaussian pulse traveling in free space normally incident upon the dielectric coating, as illustrated in Fig. 3.

We collect one time-trace at the observation point \( (P = -1.0cm) \) in the computational domain. The frequency response is obtained by using fast Fourier transformation (FFT). Moreover, the Mur’s absorbing boundary conditions (ABC) [25] is employed to truncate the open domain.

**B. Reflection coefficient**

The reflection coefficients at the observation point are calculated using the AM-DGTD method. The coating is assumed to be removed at first, that is, the whole computational domain lies in free space. The electric field data at the observation point can completely represent the incident field, \( E^{inc} \). Then, the complete model is simulated again. Because of the discontinuous interface, when the incident wave encounters the interface, a fraction of the wave energy will be reflected and part will be transmitted. Hence, the electric field data at the same observation point represents the total field \( E^{tot} \), which includes incident and reflected electric field components. The formulation for the reflected fields is given by \( E^{ref} = E^{tot} - E^{inc} \). Figure 4 illustrates the relationship of incident, total, and reflected electric field with time at the observation point.

Using FFT, the frequency responses of the incident and reflected electric fields are obtained, then the reflection coefficients \( \Gamma = E^{ref} / E^{inc} \) are achieved.

When uncertain parameters impact on the computational model, supposing \( \theta \) is independent random
parameter with probability density function (PDF). The random parameter may come from uncertain material parameters, geometrical shapes, boundary conditions, initial conditions, computational domain, etc. To model the impact of these randomness and uncertainty on the propagation of EM waves, the solution of \( E^{inc} \) and \( E^{ref} \) can be expressed as \( E^{inc}(z, t, \theta) \) and \( E^{ref}(z, t, \theta) \), they are not only a function of \((z, t)\) but also of \(\theta\). Therefore, the reflection coefficient \( \Gamma \) are uncertain and stochastic, correspondingly. The statistical moments of the solutions, such as mean and variance of \( \Gamma \) can be quantified [18]-[20] as:

\[
E[\Gamma] = \frac{1}{M} \sum_{m=1}^{M} \Gamma_m
\]

\[
Var[\Gamma] = E[(\Gamma - E[\Gamma])^2] = \frac{1}{M} \sum_{m=1}^{M} (\Gamma_m - E[\Gamma])^2,
\]

\[
Std[\Gamma] = \sqrt{Var[\Gamma]}
\]

\[
CI[\Gamma] = [E[\Gamma] - n \cdot Std[\Gamma], E[\Gamma] + n \cdot Std[\Gamma]]
\]

where \( M \) represents the samples of the stochastic parameters, which are generated by using a random number generator. \( E[\Gamma] \) stands for the mean of the random variable \( \Gamma \); \( Var[\Gamma] \) and \( Std[\Gamma] \) are the variance and standard deviation of the random variable \( \Gamma \), respectively; \( CI[\Gamma] \) denotes the confidence interval with an upper and lower bounds. The \( n \) is used as the critical value. This value is only dependent on the confidence level of the test. A typical two-sided confidence level is as \( n = 1.96 \), which corresponds to the confidence interval as 95% (see [18] for details).

C. Results and analysis

To validate the approach discussed above, the reflection coefficients are calculated by the AM-DGTD method for the sensitive variables given. The constitutive parameters are the same as in Subsection A.

Figure 5 plots the computed solution at final time \( T = 10s \) as a function of the number of adaptive elements, \( K \), and the order of the local approximation, \( N \). Comparing with analytic solutions, it shows a good agreement for different parameter sets \((N, K)\). There are two ways to improve the accuracy of \( \Gamma \): (i) keep \( N \) fixed, and increase \( K \), known as \( h \)-refinement, and (ii) keep \( K \) fixed, and increase \( N \), known as order or \( p \)-refinement [21]. As shown in the magnified image at around 1.85 GHz, simulation results using larger \( N \) or \( K \) are closer to analytic results.

Figure 6 shows the results of absolute errors (compared to analytic results). The errors decay fast with increasing \( N \) or \( K \), while both lead to a better approximation. Because of smaller errors at the extreme points, there is better manner at the maximum and minimum of the reflection coefficient \( \Gamma \), i.e., at 1.85 GHz and 3.75 GHz. But with increasing frequency, the error values become somewhat large.

Inspecting results in Fig. 6, one observes that the results of AM-DGTD scheme are clearly convergent for increasing \( K \) and/or \( N \). However, the high accuracy comes at a price. The higher accuracy needs the greater execution time. Table 1 has listed the root mean square (RMS) error and the execution time at different combinations of \((N, K)\). Consider, as an example, an error of \( O(3.0e-5) \). From the results in Table 1, we see that this can be achieved through \((2, 40), (2, 160), (3, 40), \) and \((4, 20)\) (the bold font). Comparing with these results, the combination of \((2, 40)\) has the highest order of approximation and the fastest running time in which the hosting CPU is Intel Core i3-4150 with four cores and clock speed of 3.5 GHz.

Fig. 5. The reflection coefficient for different sets of parameters \((N, K)\) compared against the exact solution.

Fig. 6. Absolute errors for different sets of parameters \((N, K)\) compared against the exact solution.
to the results of Table 1, the combination of \((N, K) = (2, 40)\) has been chosen, because of its higher accuracy and lesser time.

To analyze execution time efficiency, an ultra-thin coating is handled by the AM-DGTD method. Comparing with traditional DGTD method, the AM-DGTD approach can save more execution time. For example, when the thickness of the ultra-thin coating is 0.05 cm, the execution time is 1724.4648 and 205.0989 s, respectively, using the traditional DGTD and the AM-DGTD methods. It is clearly exhibiting that the AM-DGTD technique helps achieving very fast numerical computations.

Table 1: RMS error (GHz), total execution time (s)

<table>
<thead>
<tr>
<th>N</th>
<th>K</th>
<th>20</th>
<th>40</th>
<th>80</th>
<th>130</th>
<th>160</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>5.3615e-3</td>
<td>1.1319e-3</td>
<td>2.8409e-4</td>
<td>1.0901e-4</td>
<td>7.1152e-5</td>
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</tr>
<tr>
<td>3</td>
<td>1.8919e-3</td>
<td>2.4724e-5</td>
<td>5.1515e-5</td>
<td>1.4265e-5</td>
<td>2.7230e-5</td>
<td></td>
</tr>
<tr>
<td>5</td>
<td>5.3715e-5</td>
<td>2.7734e-5</td>
<td>1.9115e-5</td>
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<td>1.2255e-5</td>
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</tr>
<tr>
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</tr>
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<td>2.3152e-5</td>
<td>2.0652e-5</td>
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<td>1.2212e-5</td>
<td></td>
</tr>
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<td>82.5465</td>
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<td>1.3399e-5</td>
<td>1.3075e-5</td>
<td>1.5433e-5</td>
<td>1.3684e-5</td>
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</tr>
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</tr>
</tbody>
</table>

IV. AM-DGTD FOR STOCHASTIC PARAMETERS

The actual form of flaws in the dielectric coating may not be known, causing the introduction of the uncertainty in some sensitive parameters. These uncertainties can be based on pure speculation, on measured data, or on other available information. In this section, a few typical uncertain flaws are discussed and validated by the AM-DGTD approach. To reduce the complexity of the simulations, the uncertain flaws are expressed as stochastic variables, characterized with appropriate probability distributions. And these stochastic variables are statistically independent.

A. Uncertainty in material

Deterministic solutions require accurate input parameters, however, there always exist uncertain material properties such as from imprecise measurement or manufacturing. In this section, the case of uncertainty in permittivity is considered, that is, tolerances and uncertainties lie in the relative permittivity \(\varepsilon_r\).

For the model in Fig. 3, a randomness in the relative permittivity of the dielectric coating is assumed as [9]:

\[
\varepsilon_r(x, \theta) = \begin{cases} 
1 & x \notin \text{coating} \\
4 \cdot (1 + \frac{0.1 \cdot \theta^2}{1 + \theta^2}) & \text{otherwise}
\end{cases}, \quad (9)
\]

where \(\theta\) is a Gaussian variable with zero mean and unit variance. With the uncertainty of the formula (9), the relative permittivity \(\varepsilon_r\) is guaranteed to remain positive in the domain of the dielectric coating. Note that our concern is not the correctness of the probabilistic law chosen in (9) for the uncertainty of the \(\varepsilon_r\), since for any reasonable law, the techniques presented in this paper should work equally well.

Other parameters are fixed as the computational domain \(\Omega = [-6.0 \text{cm}, 6.0 \text{cm}]\), the thickness of the coating \(d = 2.0 \text{cm}\), the position of the observed point \(P = -0.1 \text{cm}\), the total computational time \(T = 10s\).

To study the relationship of the deviation of reflection coefficient \(\Gamma\) and permittivity \(\varepsilon_r\), another 1000 independent random variables are used. The results of uncertainty quantification of reflection coefficient with different deviations in permittivity \(\varepsilon_r\) over a 0-4 GHz frequency range are shown in Fig. 7. It is obvious that the minimum
reflection point, i.e., 3.75 GHz, turn worse with the tolerances of $\varepsilon_r$ increase. When the tolerance of $\varepsilon_r$ is changed, the value of this point is increased, i.e., the value of $\varepsilon_r$ decreases, this point shifts a bit to higher frequency; when the values of $\varepsilon_r$ increases, the point shifts a bit to lower frequency. The increased offset indicates an increased tolerance existing in the results.

Table 2: Uncertainty quantification of $\Gamma$ versus different deviation in $\varepsilon_r$

<table>
<thead>
<tr>
<th>$\Delta\varepsilon_r \Gamma$</th>
<th>Min</th>
<th>Max</th>
<th>Mean</th>
<th>95% CI</th>
</tr>
</thead>
<tbody>
<tr>
<td>-0.4056</td>
<td>0.4737</td>
<td>0.5999</td>
<td>0.5595</td>
<td>[0.4905, 0.6286]</td>
</tr>
<tr>
<td>-0.2831</td>
<td>0.5287</td>
<td>0.5999</td>
<td>0.5735</td>
<td>[0.5311, 0.6159]</td>
</tr>
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<td>-0.1388</td>
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<tr>
<td>0</td>
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<td>0.1287</td>
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<td>0.6267</td>
<td>0.6098</td>
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<tr>
<td>0.2994</td>
<td>0.5999</td>
<td>0.6474</td>
<td>0.6214</td>
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</tr>
<tr>
<td>0.4232</td>
<td>0.5999</td>
<td>0.6668</td>
<td>0.6286</td>
<td>[0.5882, 0.6690]</td>
</tr>
</tbody>
</table>

Fig. 7. Distribution of $\Gamma$ with respect to different deviation in $\varepsilon_r$.

B. Uncertainty in thickness of coating

Suppose that the physical thickness of the dielectric coating is unknown, i.e., when there is uncertainty associated with the position of the right boundary of the coating. Uncertainty in the thickness of $d$ can be modeled by choosing $x^{right}$ a random variable and keeping $x^{left}$ as constant. And the randomness $x^{right}$ is a random parameter with some associated PDF.

The uncertain deviation in the right boundary position is chosen, so that the thickness of the coating is of variable value. The position of the right boundary is assumed to be $x^{right} = d + g(\theta)$, this allows $x^{right}$ to be positioned on both sides of the mean position. Here $g(\theta)$ (the deviation in thickness $\Delta d$) is a uniform variable. The mean is chosen as $\pm 0.1$, $\pm 0.2$ and $\pm 0.3$ with a tolerance of $\pm 0.01$, i.e., $g(\theta)/\Delta d$ is uniformly distributed in the interval $[\pm 0.09cm, \pm 0.11cm]$, $[\pm 0.19cm, \pm 0.21cm]$, and $[\pm 0.29cm, \pm 0.31cm]$, respectively.

Figure 8 shows how the uncertain deviation in thickness of the dielectric coating affects the reflection coefficient $\Gamma$ at the observation point $P = -0.1cm$, which are computed using the AM-DGTD formulation outlined in Section III. In the frequency 0-5 GHz range, the magnitude of the mean of $\Gamma$ for 1000 samples of the random thickness is shown in the Fig. 8. When $\Delta d = 0$, the numerical results are in good agreement with analytic solutions. With the thickness $d$ decreasing, e.g., the average thickness is from 2.3cm to 1.7cm, it has been observed obviously that the curves of the reflection coefficient $\Gamma$ are extended towards the $x$ axis and the minimum reflection frequency points shift from low to high frequency.

C. Random holes in coating

In this experiment, the AM-DGTD method is employed to analyze the reflection coefficients of imperfect dielectric coating caused by blotch, bubble, or recessed hole on a film. The different number, different radius, and random positions of holes in the dielectric coating of the range of $[0.0cm, 2.0cm]$ are simulated.

1) Single hole

The hole parameters are assumed as $\varepsilon_h = 1$, $\mu_h = 1$. The arbitrary mesh is represented in Fig. 9, the radius and random position of the hole are statistically independent. The number of the elements is set as $K = 40$. Degrees of two ($N = 2$) polynomials are enough to ensure that convergence is achieved in the physical space. The randomness of the holes can be incorporated into the AM-DGTD method by introducing uncertainty in the local computational mesh. To accomplish this, a displacement mesh is employed around each uncertain portion of the geometry.
Three cases with random position and radius $r$ as 0.025cm, 0.05cm, and 0.1cm are considered, respectively. Comparisons of the mean of the reflection coefficient $\Gamma$ with and without these random holes over a 0-10 GHz frequency range are presented in Fig. 10. When the radius $r$ is increased, the mean $\Gamma$ becomes worse. For instance, the maximum reflection coefficient at 1.85 GHz is decreased, and the minimum value at 3.75 GHz is increased. And the curves of frequency-domain response are shifted a bit to higher frequency.

2) Multiple holes

An arbitrary mesh with two uncertain holes is represented in Fig. 11, all parameters are same as that of single-hole case. With the increasing of hole radius, the mean of reflection coefficient $\Gamma$ becomes worse. For example, the value of the maximum becomes smaller and the value of the minimum becomes bigger.

The properties of results are similar to those of the single-hole case, but for the two random holes, the curve shifts more bit to the higher frequency, as shown in Fig. 12. The comparison results of the mean of the reflection coefficient of the model with 0-3 random holes are summarized in Fig. 13. It has been found that the mean $\Gamma$ becomes worse with the number of the holes increasing.

Figure 14 shows the estimation of mean $\Gamma$ and 95% CI for the imperfect dielectric coating with random number of holes computed by the AM-DGTD method. And the radius of hole is supposed to equal to 0.1cm, and the positions of hole are randomly selected. The limits of the 95% CI are obtained from the PDF, such that...
the 2.5% of the data goes beyond the lower and upper bounds, respectively. The mean $\Gamma$ of 2.5% quartile near the minimum is observed to perturb slightly and the mean $\Gamma$ of 97.5% near the maximum is observed to shift slightly, all indicating an increased uncertainty exists in the results.

$$\text{Fig. 14. The deviation in the mean and in 95\% confidence intervals for } \Gamma \text{ when the solution with stochastic holes is compared against the solution without holes: (a) one hole, (b) two holes, and (c) three holes.}$$

Figure 15 (a) shows the mean of the $\Gamma$ for the dielectric coating without holes by using the AM-DGTD method. The mean $\Gamma$ with only one uncertain hole at the maximum 1.85 GHz are concentrated in the interval [0.581, 0.598], however, the values drop to range of [0.562, 0.597] and [0.544, 0.596] for the two and three holes, respectively, as shown in Figs. 15 (b)-(d). These numerical results are agreed with the physical explanation. In term of the value of the interval, we clearly see the smaller mean $\Gamma$ have worse flaws corresponding to a larger number of holes.

From Fig. 16, it is clearly seen that the mean $\Gamma$ at different frequencies is different for different numbers of random holes. For $0 \leq \Gamma \leq 0.4$, the values are increased with increasing of the number of the random hole; the $\Gamma = 0.5$ results slightly changed; and for $0.5 < \Gamma \leq 0.6$, the values are decreased with the increasing of the number of the random hole.

$$\text{Fig. 15. Distribution of the reflection coefficient with different number of random holes: (a) the analytic solution of } \Gamma \text{ without hole, (b) the distribution of } \Gamma \text{ with one random hole at 1.85 GHz, (c) the distribution of } \Gamma \text{ with two holes, and (d) the distribution of } \Gamma \text{ with three random holes.}$$

$$\text{Fig. 16. The relationship between the mean of } \Gamma \text{ and the number of hole.}$$

V. CONCLUSION

In this paper, we discussed the use of an adaptive mesh integrated into discontinuous Galerkin method to research the impact of uncertain flaws in the 1D propagation problems. These sources of uncertain flaws are considered: the varied relative permittivity, thickness of coating and multivariable holes in the medium. The adaptive spatial discrete algorithm for the high order accurate DGTD method has been developed to solve
these irregular, complex, and random flaws in the dielectric coating.

The adaptive mesh technology, which generated by the displacement of nodes and the insertion of nodes, can provide efficient and optimum number of mesh elements as compared to simple uniform mesh. The simulation results have been shown that the new approach can save computational resources because of avoiding redundant division of the computational domain. Three typical experiments validate significant advantages of the AM-DGTD approach and show potential for further study of uncertain problems.

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