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Uncertainty Analyses in the Finite-Difference Time-Domain Method

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Abstract—Providing estimates of the uncertainty in results obtained by Computational Electromagnetic (CEM) simulations is essential when determining the acceptability of the results. The Monte Carlo method (MCM) has been previously used to quantify the uncertainty in CEM simulations. Other computationally efficient methods have been investigated more recently, such as the polynomial chaos method (PCM) and the method of moments (MoM). This paper introduces a novel implementation of the PCM and the MoM into the finite-difference time-domain method. The PCM and the MoM are found to be computationally more efficient than the MCM, but can provide poorer estimates of the uncertainty in resonant electromagnetic compatibility data.

Index Terms—Computational electromagnetism, finite-difference time domain (FDTD), method of moments (MoM), Monte Carlo, polynomial chaos, uncertainty analysis.

I. INTRODUCTION

COMPUTATIONAL electromagnetic (CEM) simulations rely on sets of input parameters, which often have an associated uncertainty. These uncertainties may arise from a lack of precise knowledge of the material parameters or geometries that are being modeled. Uncertainties in these input parameters lead to uncertainties in the output of the CEM simulations. This type of uncertainty is often known as parameter uncertainty. In this paper, a determination of the parameter uncertainty in the results of finite-difference time-domain (FDTD) simulations will be made. Quantifying the uncertainty in the output of interest amounts to quantifying the standard deviation of the output. Uncertainty analyses provide the quantitative level of confidence that may be held in the results of CEM simulations. This information is essential when determining whether the results are acceptable or useful.

Previous research has already highlighted the importance of quantifying uncertainty in CEM [1]–[4]. This research uses the Monte Carlo method (MCM), which is generally accepted as being an accurate uncertainty analysis (UA) method, to test the performance of other computationally efficient UA methods. Chauvière published work involving the implementation of the polynomial chaos method (PCM) into a higher order discontinuous Galerkin solution of Maxwell’s equations [1]. The PCM was found to accurately quantify the output uncertainty, while being more computationally efficient than the MCM. Chauvière’s work, however, only estimated the output uncertainty due to one uncertain input parameter. The accuracy and computational efficiency of this method with increased numbers of uncertain input parameters needs to be analyzed.

More recently, Ajayi has discussed the use of a direct solution technique (DST) to quantify uncertainty [2]. This technique applies the probabilistic method of moments (MoM) [5], [6] to different CEM schemes, such as the transmission line matrix (TLM) method. Ajayi used the DST to estimate the uncertainty in the frequency of the first resonance for simple electromagnetic problems [2]. The DST was found to work well for small parameter variations, giving results that are in agreement with results obtained from the MCM [2].

This paper outlines novel implementations of the PCM and the MoM into the FDTD method. It is possible to implement these statistical methods into other CEM techniques. The uncertainty in the output of simulations performed using different CEM methods and different levels of accuracy will be of a similar size. However, with the results of different simulations formed by different CEM techniques, the uncertainty in the output will have a dependence on the method used and the accuracy with which the simulation is performed. This paper considers only the FDTD method so that a fair analysis of the PCM, the MoM, and the MCM can be formed.

In this paper, the UA methods are used to obtain the uncertainty in the output electric field viewed in the frequency domain. In the first of two examples, the UA methods are used to estimate the uncertainty in the electric field that penetrates a shielded enclosure containing a printed circuit board (PCB), at around 1.8 GHz. This example is fairly simple, having only one uncertain input parameter and encompassing only a few resonant features. The second electrically large example considers three uncertain inputs, encompassing many more resonant modes. The PCM and the MoM are compared to the MCM in terms of their ability to accurately quantify the uncertainty and their computational expense. The uncertain input parameters considered in this paper are all assumed to be uncorrelated. When considering correlated variables, United Kingdom Accreditation Service (UKAS) [6] suggests grouping all correlated variables into one single grouped variable, which will subsequently be uncorrelated with all other uncertain variables in the sample.

In this paper, the uncertainty in the output of interest is a frequency response curve. The feature selective validation (FSV) method [7], [8] is used in this paper to determine the similarity of the uncertainty curves formed from the different UA methods. This method is a numerical technique, which determines how similar two curves are in terms of their amplitude and feature differences. The amplitude and feature differences between the curves are combined to give a general difference measure (GDM). This method currently forms part of a draft standard for the verification and validation of CEM models [9].
II. UA METHODS

A. Monte Carlo Method

To determine the uncertainty in an FDTD simulation via the MCM, the probability density functions (PDFs) associated with the uncertain input parameters must first be sampled many times. In this paper, latin hypercube sampling (LHS) [10] is chosen as the preferred sampling method. This has been shown to produce a converged solution more quickly than other sampling methods [10]. The samples, obtained from the PDFs, form sets of input parameter values: one FDTD simulation is performed for each set. The outputs formed from each simulation are combined to form the output mean and standard deviation. The uncertainty in the output is represented by the standard deviation.

It is well known that the MCM has slow convergence, and as such, it is a computationally expensive method. The mean and uncertainty, formed using the MCM, converge for large numbers of samples. Once convergence has been reached, the use of more samples does not change the mean and uncertainty significantly. In this paper, the FSV method is used, in a novel way, to determine when the MCM has reached convergence. The mean and uncertainty formed in this paper are frequency response curves. The FSV method is used to compare the respective mean and uncertainty frequency response curves after every $N$ simulations. Convergence is reached when the mean and uncertainty curves are determined to be “excellent” matches to the respective mean and uncertainty curves produced after $N$ simulations previously. In the examples in this paper, $N = 50$ and convergence is reached when the comparisons of the respective mean and uncertainty curves produce a GDM < 1.5. This novel way of using FSV provides an accurate, consistent, and impartial way of determining convergence, when the output of interest is a curve.

B. Polynomial Chaos Method

The concept of homogeneous chaos was first introduced by Wiener [11]. Homogeneous chaos uses Hermite polynomials to represent stochastic processes that depend on uncertain input parameters, which follow normal distributions [12]. The PCM is a more generalized method for dealing with inputs that are not necessarily normally distributed. In the PCM, certain orthogonal basis polynomials are selected, depending on the distributions of the random input variables [12]. A function depending on the uncertain random variables can be expanded in terms of the selected polynomials. This expansion, which is known as the Wiener–Askey chaotic expansion [12], casts the uncertainty in the output into the orthogonal polynomials alone. The orthogonality of the polynomial basis set can reduce stochastic differential equations to a set of deterministic differential equations that can be solved numerically [12].

Xiu and Karniadakis [12] found the PCM to be computationally cheaper than the MCM. However, they noted that the method’s efficiency is problem specific [13].

1) Wiener–Askey Chaos: Any second-order random process $X$, depending on some random event $\theta$, can be represented as [12]

$$X(\theta) = \sum_{i=0}^{\infty} c_i \psi_i(\zeta(\theta))$$  \hspace{1cm} (1)

where $c_i$ are constant coefficients and

$$\zeta(\theta) = (\zeta_1(\theta), \zeta_2(\theta), \ldots)$$  \hspace{1cm} (2)

represents a vector containing an infinite number of independent random variables [12]. The polynomial basis sets $\{ \psi_i \}$, corresponding to the random variables $\zeta(\theta)$, are chosen from the Askey-scheme, which can be found in [12].

The polynomial basis sets are all orthogonal, which implies that

$$\langle \psi_i \psi_j \rangle = \langle \psi_i^2 \rangle \delta_{ij}$$  \hspace{1cm} (3)

where $\delta_{ij}$ is the Kronecker delta and the inner product $\langle ., . \rangle$ is defined as [12]

$$\langle f(\zeta) g(\zeta) \rangle = \int f(\zeta) g(\zeta) w(\zeta) d\zeta.$$  \hspace{1cm} (4)

The weighting function $w(\zeta)$ corresponds to the choice of polynomial basis $\{ \psi_j \}$.

2) General Polynomial Chaos: The Weiner–Askey chaotic expansion (1) can be used to solve stochastic differential equations [12]. Let $u(x, t, \theta)$ be a solution of the stochastic differential equation

$$\mathcal{L}(x, t, \theta) u(x, t, \theta) = f(x, t, \theta)$$  \hspace{1cm} (5)

where $x$, $t$, and $\theta$ represent position, time and some random event, respectively. The symbol $\mathcal{L}$ represents some differential operator and $f$ is a source term [12]. The solution $u$ may be regarded as a random process and expanded as [12]

$$u(x, t, \theta) = \sum_{i=0}^{P} u_i(x, t) \psi_i(\zeta(\theta)).$$  \hspace{1cm} (6)

For practical applications, the infinite sum in (1) has been truncated at P here. If $d$ is the order of the highest order polynomial used in the expansion and $n$ is the dimension of the random variable $\zeta$, then

$$P + 1 = \frac{(n + d)!}{n!d!}.$$ \hspace{1cm} (7)

Substituting the expansion of $u$ into (5) yields

$$\mathcal{L}(x, t, \theta) \sum_{i=0}^{P} u_i(x, t) \psi_i(\zeta(\theta)) = f(x, t, \theta).$$  \hspace{1cm} (8)

The inner product of both sides of this equation can be formed with $\psi_k$ to give

$$\langle \mathcal{L}(x, t, \theta) \sum_{i=0}^{P} u_i(x, t) \psi_i(\zeta(\theta)), \psi_k \rangle = \langle f(x, t, \theta), \psi_k \rangle.$$  \hspace{1cm} (9)

The orthogonality of the basis polynomials reduces (9) to a set of $(P + 1)$ differential equations: one for each $u_i$ [12]. These equations are deterministic [12] and can, therefore, be solved
numerically. Once each \( u_i \) is found, the mean and variance of \( u(x, t, \theta) \) can be calculated. The mean is calculated as [1]

\[
\bar{u}(x, t, \theta) = \langle u(x, t, \theta), 1 \rangle = \sum_{i=0}^{P} u_i \langle \psi_i, 1 \rangle
\]

\[
= \sum_{i=0}^{P} u_i \langle \psi_i, \psi_0 \rangle = \sum_{i=0}^{P} u_i \delta_{i0} = u_0
\]

using the fact that \( \psi_0 = 1 \) for all polynomial bases. The variance can be obtained in a similar way by first calculating

\[
\langle u(x, t, \theta), u(x, t, \theta) \rangle = \sum_{i=0}^{P} \sum_{j=0}^{P} u_i u_j \langle \psi_i, \psi_j \rangle
\]

\[
= \sum_{i=0}^{P} \sum_{j=0}^{P} u_i u_j \delta_{ij} \langle \psi_i^2 \rangle = \sum_{i=0}^{P} u_i^2 \langle \psi_i^2 \rangle.
\]

(11)

The variance of \( u(x, t, \theta) \) is therefore [1]

\[
\sigma^2 = \langle u(x, t, \theta), u(x, t, \theta) \rangle - \langle u(x, t, \theta), 1 \rangle^2
\]

\[
= \sum_{i=0}^{P} u_i^2 \langle \psi_i^2 \rangle - u_0^2 = \sum_{i=0}^{P} u_i^2 \langle \psi_i^2 \rangle.
\]

(12)

The uncertainty in \( u(x, t, \theta) \) is the standard deviation \( \sigma \).

3) General Polynomial Chaos in FDTD: General polynomial chaos has already been successfully applied to computational fluid dynamics (CFD) [14, 15] and specific areas of CEMs [1]. The application of general polynomial chaos to 1-D FDTD is given here, and these arguments can be easily generalized to three dimensions.

Maxwell’s equations for a wave propagating in a linear isotropic homogeneous material along the \( x \)-axis in 1-D are

\[
\frac{\partial E_y}{\partial t} = -\frac{1}{\mu} \frac{\partial H_z}{\partial x}
\]

\[
\frac{\partial H_z}{\partial t} = -\frac{1}{\epsilon} \left( \frac{\partial E_y}{\partial x} + \sigma E_y \right).
\]

As usual, \( H_z(x, t) \) represents the magnetic field oriented in the \( z \)-direction, at a position \( x \) and time \( t \). Similarly, \( E_y \) represents the electric field oriented in the \( y \)-direction. The symbols \( \mu \), \( \epsilon \), and \( \sigma \) represent the permeability, permittivity, and conductivity of the medium in which the electromagnetic fields propagate. CEM models seek to solve these two coupled equations to find approximations for \( H_z(x, t) \) and \( E_y(x, t) \). If there are uncertain input parameters, then the solutions will depend on some random event \( \theta \). The uncertain field solutions can, therefore, be represented as \( H_z(x, t, \theta) \) and \( E_y(x, t, \theta) \). The solutions to Maxwell’s equations may be found by using the FDTD scheme, first proposed by Yee [16].

In Yee’s scheme, the temporal and spatial partial derivatives in Maxwell’s equations are approximated using central difference approximations. The problem space is discretized into cells of length \( \Delta x \), and the time is split into discrete intervals \( \Delta t \). This yields two update equations, which form the basis of the 1-D FDTD solution. The update equations formed are

\[
H_z^{n+1/(2)}(j + 1/2, \theta) = H_z^{n-(1/2)}(j + 1/2, \theta) - \gamma [E_y^n(j + 1, \theta) - E_y^n(j, \theta)]
\]

(15)

and

\[
E_y^{n+1}(j, \theta) = \alpha E_y^n(j, \theta) + \beta \left[ H_z^{n+1/(2)}(j - 1/2, \theta) - H_z^{n+1/(2)}(j + 1/2, \theta) \right].
\]

(16)

The shorthand notation \( H_z^n(j, \theta) = H_z(j \Delta x, n \Delta t, \theta) \) is used in the previous equations, where \( j \) and \( n \) are positive integers. The material properties of the medium in which the fields propagate are represented by

\[
\alpha = \alpha(x, \theta) = \frac{1 - \sigma(x, \theta) \Delta t/2 \epsilon(x, \theta)}{1 + \sigma(x, \theta) \Delta t/2 \epsilon(x, \theta)}
\]

(17)

\[
\beta = \beta(x, \theta) = \frac{\Delta t}{\Delta x \mu(x, \theta)}
\]

(18)

and

\[
\gamma = \gamma(x, \theta) = \frac{\Delta t}{\Delta x \mu(x, \theta)}
\]

(19)

The material properties \( \mu, \epsilon, \) and \( \sigma \) all depend on the uncertain parameter \( \theta \). This dependence is defined by the PDFs of the input parameters.

The 1-D FDTD update equations defined above are used to obtain solutions for the electric and magnetic fields, subject to some electric field source. The field solutions can be expanded in terms of the appropriate orthogonal polynomials \( \{ \psi_i \} \) to separate the dependence of the field on the random parameter \( \theta \) from the dependence on time and position. The field solutions are expanded as

\[
E_y^n(j, \theta) = \sum_{i=0}^{P} e_i^n(j) \psi_i(\zeta(\theta))
\]

(20)

and

\[
H_z^n(j, \theta) = \sum_{i=0}^{P} h_i^n(j) \psi_i(\zeta(\theta)).
\]

(21)

The coefficients \( e_i^n(j) \) and \( h_i^n(j) \) must be found so that the mean and uncertainty of the output fields can be formed. The following discussion outlines how to obtain the field coefficients \( e_i^n(j) \) and \( h_i^n(j) \).

Expansions (20) and (21) can be substituted into (15) and (16) to obtain

\[
\sum_{i=0}^{P} h_i^n(1/2)(j + 1/2) \psi_i(\zeta(\theta)) = \sum_{i=0}^{P} \left[ h_i^{n-(1/2)}(j + 1/2) \right]
\]

\[
\times \psi_i(\zeta(\theta)) + \gamma (e_i^n(j + 1) - e_i^n(j)) \psi_i(\zeta(\theta))
\]

(22)
and
\[ \sum_{i=0}^{P} e_{i}^{n+1}(j) \psi_{i}(\theta) = \sum_{i=0}^{P} [\alpha e_{i}^{n}(j) \psi_{i}(\theta) + \beta \left( h_{i}^{n+1/2}(j - 1/2) - h_{i}^{n+1/2}(j + 1/2) \right) \psi_{i}(\theta)]. \] (23)

At this point, the electric field source \( E_{s} \) may be added into (23). If this field source has some associated uncertainty, then it may be expanded as
\[ E_{s}^{n}(j, \theta) = \sum_{i=0}^{P} e_{i,s}^{n}(j) \psi_{i}(\theta). \] (24)

Adding this source term into (23) yields
\[ \sum_{i=0}^{P} e_{i}^{n+1}(j) \psi_{i} = \sum_{i=0}^{P} \left[ \alpha e_{i}^{n}(j) \psi_{i} + e_{i,s}^{n}(j) \psi_{i} + \beta \left( h_{i}^{n+1/2}(j - 1/2) - h_{i}^{n+1/2}(j + 1/2) \right) \psi_{i} \right]. \] (25)

Taking the inner product of both sides of (22) and (25) with some test polynomial \( \psi_{k} \), where \( k \) is an integer in the range \( 0 \leq k \leq P \), reduces the equations to
\[ h_{k}^{n+1/2}(j + 1/2) = h_{k}^{n-1/2}(j + 1/2) - \frac{1}{\langle \psi_{k}^{2} \rangle} \sum_{i=0}^{P} \left[ \left( e_{i}^{n}(j + 1) - e_{i}^{n}(j) \right) \langle \gamma \psi_{i} \psi_{k} \rangle \right] \] (26)
and
\[ e_{k}^{n+1}(j) = e_{k,s}^{n+1}(j) + \frac{1}{\langle \psi_{k}^{2} \rangle} \sum_{i=0}^{P} \left[ \left( e_{i}^{n}(j) \langle \alpha \psi_{i} \psi_{k} \rangle - \langle \alpha \psi_{i} \psi_{k} \rangle \right) e_{i}^{n+1}(j) \right] + \left( h_{k}^{n+1/2}(j - 1/2) - h_{k}^{n+1/2}(j + 1/2) \right) \langle \beta \psi_{i} \psi_{k} \rangle. \] (27)

The previous two equations make use of the orthogonality relation set out in (3). The material properties \( \alpha \), \( \beta \), and \( \gamma \) remain in the inner products due to their dependence on \( \theta \).

Equations (26) and (27) can be used to calculate \( e_{k}^{n+1}(j) \) and \( h_{k}^{n+1/2}(j + 1/2) \) for all \( n \) and \( j \), and \( k = 0, \ldots, P \). In order to do this, the inner products \( \langle \alpha \psi_{i} \psi_{k} \rangle \), \( \langle \beta \psi_{i} \psi_{k} \rangle \), and \( \langle \gamma \psi_{i} \psi_{k} \rangle \) must first be calculated for all \( i, k = 0, \ldots, P \). These can be calculated using numerical integrations as a preprocessing; the update equations can then be used in a similar manner to the leapfrog scheme used in conventional FDTD. The source term must also be calculated at each time step as
\[ e_{k,s}^{n}(j) = \frac{1}{\langle \psi_{k}^{2} \rangle} \langle E_{s}^{n}(j, \theta) \psi_{k}(\theta) \rangle. \] (28)

To complete the update, update equations are required at the boundary. The examples in this paper consider the scattering of electromagnetic fields in free space, and therefore, the first-order Mur absorbing boundary condition [17] is chosen as a relevant boundary condition. At the lower boundary (where \( j = 0 \), the usual Mur update equation is
\[ E_{y}^{n+1}(0) = E_{y}^{n}(1) + \sqrt{\frac{\beta(1, \gamma)(1) - 1}{\beta(1, \gamma)(1) + 1}} \left( E_{y}^{n+1}(1) - E_{y}^{n}(0) \right). \] (29)

Introducing uncertainty into this equation via dependence on \( \theta \) yields
\[ E_{y}^{n+1}(0, \theta) = E_{y}^{n}(1, \theta) + \sqrt{\frac{\beta(1, \gamma)(1, \theta) - 1}{\beta(1, \gamma)(1, \theta) + 1}} \left( E_{y}^{n+1}(1, \theta) - E_{y}^{n}(0, \theta) \right). \] (30)

As before, the field terms can be expanded using the chaotic expansion, and an inner product of both sides of the resulting equation can be taken with \( \psi_{k} \). Carrying out these two steps gives
\[ e_{k}^{n+1}(0) = e_{k}^{n}(1) + \frac{1}{\langle \psi_{k}^{2} \rangle} \sum_{i=0}^{P} \left( e_{i}^{n+1}(1) - e_{i}^{n}(0) \right) \langle \xi(1, \theta) \psi_{i} \psi_{k} \rangle \] (31)
where \( \xi \) is defined as
\[ \xi(1, \theta) = \sqrt{\frac{\beta(1, \gamma)(1, \theta) - 1}{\beta(1, \gamma)(1, \theta) + 1}}. \] (32)

Usually, at the boundaries, the properties of the medium are those of free space: there is no uncertainty in the parameter values near the boundary. Thus, \( \xi \) has no \( \theta \) dependence and \( \langle \psi_{i} \psi_{k} \rangle = \delta_{ik} \langle \psi_{k}^{2} \rangle \). The update equation for the electric field at the lower boundary reduces to
\[ e_{k}^{n+1}(0) = e_{k}^{n}(1) + \xi(1) \left( e_{k}^{n+1}(1) - e_{k}^{n}(0) \right) \] (33)
for \( k = 0, \ldots, P \). Using a similar argument, the electric field update equation at the upper boundary (\( j = N \)) is
\[ e_{k}^{n+1}(N) = e_{k}^{n}(N - 1) + \xi(N - 1) \left( e_{k}^{n+1}(N - 1) - e_{k}^{n}(N) \right). \] (34)

The coefficients \( e_{k}^{n}(j) \) may now be calculated using (26), (27), (33), and (34). Similar equations to (10) and (12) are formed for the mean value of \( E_{y}^{n}(j, \theta) \)
\[ \bar{E}_{y} = e_{0}^{n}(j) \] (35)
and the variance
\[ \sigma^{2} = \sum_{i=1}^{P} \left( e_{i}^{n}(j) \right)^{2} \langle \psi_{i}^{2} \rangle. \] (36)

The PCM is easily applied to FDTD because the update equations follow a very similar form to those in conventional FDTD. The only extra step comes in calculating the integrals that correspond to the inner products in the update equations (26) and (27). The application of the PCM to FDTD given here may be easily generalized to 3-D.

One problem with the PCM arises when trying to form the mean and uncertainty of some related quantity, which does not depend linearly on \( E_{y}^{n}(j, \theta) \). For example the output of interest.
may be the absolute value of $E_y^n(j, \theta)$. The mean $\mu$ of this output is formed as

$$\mu = \langle |E_y^n(j, \theta)|, 1 \rangle = \left( \sum_{j=0}^{P} |e_y^n(j)\psi_j(\chi)\rangle |\psi_0\rangle \right). \quad (37)$$

The absolute value within the inner product prevents the orthogonality of the basis polynomials from being used to form a simple relationship for the mean. The mean will have to be calculated using a numerical integration over the uncertain parameter space $\theta$. This presents a problem with the PCM: the calculation of the mean of the output of interest is not always trivial. Some of the mathematical simplicity of the PCM has been lost by trying to form the mean of the quantity $|E_y^n(j, \theta)|$. The mean must be calculated using a numerical integration at each frequency point, which requires extra computational time. This extra computational expense is, however, small compared to the PCM simulation runtime. Once this mean has been calculated, the variance $\sigma^2$ may be calculated using the standard definition

$$\sigma^2(|E_y^n(j, \theta)|) = \langle (|E_y^n(j, \theta)| - \mu)^2 \rangle. \quad (38)$$

Further numerical integrations are required to obtain this variance. The uncertainty in $|E_y^n(j, \theta)|$ is the square root of this variance.

The PCM performs one large simulation, storing a factor $(P + 1)$ more field coefficients and $(P + 1)^2$ more material parameter values (via the inner products) than the MoM and the MCM. If there are, for example, three uncertain inputs and the Wiener–Askey chaos expansion is truncated at first order, then $P = 3$, and the PCM will require around 16 times more memory than the MoM and the MCM to store the material inner products.

The computational time required by the PCM is somewhat more difficult to analyze. Since there are $(P + 1)$ field coefficients to solve for and each calculation requires the sum of the product of $(P + 1)$ field coefficients with $(P + 1)$ material inner products [see (26) and (27)], the computational time required by the PCM will be approximately $(P + 1)^2$ greater than that of a single FDTD simulation. In addition to this, extra computational time is required to calculate the material inner products at each point in the problem space and the output variance from the field coefficients. These extra numerical integrations add to the overall computational effort of the PCM; however the computational time will scale approximately with $(P + 1)^2$. This computational time will be greater than that required for the MoM, but should be less than that required for the MCM.

C. Method of Moments

The MoM is another approximate UA method, which is similar to the method outlined in [6] for the determination of uncertainty in practical electromagnetic compatibility (EMC) measurements. It is the internationally accepted method outlined in [5] for the propagation of uncertainties through a model.

The MoM uses a first-order Taylor series expansion of the output electric field $E_y$ about the mean input parameter values

$$E_y(p_1, \ldots, p_n) = c_1 p_1 + \cdots + c_n p_n \quad (39)$$

where $c_i$ represent the sensitivity derivatives of each parameter $p_i$ evaluated at the mean parameter values $\bar{p}_i$. For the purposes of this discussion, it is assumed that there are $n$ input parameters for the FDTD simulation.

The mean output electric field $\bar{E}_y$ is calculated by performing one simulation with all input parameters taking on their mean values [5], [6]. To calculate the sensitivity derivative $c_i$, in (39), an FDTD simulation must be performed with all parameters taking on their mean values, except for the parameter $p_i$. In this simulation, the parameter $p_i$ is perturbed slightly from its mean value to give $p_i = \bar{p}_i + \Delta$. This simulation will produce a perturbed output electric field $\tilde{E}_y$. The sensitivity derivative $c_i$ is calculated using the finite-difference approximation

$$c_i = \frac{\tilde{E}_y - \bar{E}_y}{\Delta}. \quad (40)$$

If $u_i$ is the uncertainty in the parameter $p_i$, determined from the PDF of $p_i$, then the uncertainty in the mean output of the FDTD simulation is calculated as [5], [6]

$$\sigma = \left( \sum_{i=1}^{n} c_i^2 u_i^2 \right)^{1/2}. \quad (41)$$

The accuracy of the uncertainty estimate relies on the relationship between the uncertain inputs and the output of interest being linear. The method is also dependent on the size of the perturbation $\Delta$ that is used. It has been previously suggested that using a perturbation $\Delta = u_i$ is appropriate [5], [6].

The MoM requires the same amount of computational memory as the MCM. For a simulation with $n$ uncertain input parameters, the MoM requires $n + 1$ FDTD simulations.

In the next section, the UA methods described earlier are used to determine the mean and uncertainty in the output of an EMC example. The output of interest in the following examples is the frequency response of the normalized electric field. This normalized field is formed by taking the ratio of the absolute value of the specified electric field to the input excitation in the frequency domain. The FDTD simulations are all performed with $100 \times 100 \times 100$ cells and 10 000 time steps.

III. Example 1: An EMC Example

Fig. 1 shows the setup of the EMC example. An electric field is excited from a dipole (oriented in the $z$-direction) and scattered off a shielded enclosure, with an aperture in the front face. To describe the coordinates of the shielded enclosure and the aperture, the origin is assigned to the bottom right-hand corner of the front face, which itself is in the $y-z$ plane. The width of the box is $y = 38$ cm, the depth is $x = 40$ cm, and the height is $z = 15$ cm. For the purposes of the following discussion, the points in this coordinate system have units of centimeters, referenced from the origin. The bottom right-hand corner of the aperture is at the point $(0, 19, 5)$ and extends by a width $a_y = 15$ cm and a height $a_z = 4$ cm. Using this coordinate
system, the center of the dipole is at ($-20, 26, 7$), this is 20 cm away (in the $x$-direction) from the center of the aperture. The arms of the dipole are each 7 cm in length, with a radius of 1 mm. The voltage source at the center of the dipole has an amplitude of $V_0 = 2$ V over a load of 50 $\Omega$. The input excitation is a Gaussian of the form

$$V = V_0 \exp \left(-\frac{4 \ln 2 (t - t_0)^2}{f w h h^2}\right) \quad (42)$$

where $t_0 = 6.67 \times 10^{-10}$ s is the onset time and $f w h h = 2.78 \times 10^{-10}$ s is the full width of the Gaussian pulse at half the height of the maximum amplitude.

The enclosure represents the shielding exterior of a typical electronic system containing a PCB. Using the coordinate system outlined before, the PCB is oriented in the $x$-$z$ plane, extending from the point (5, 14, 2) a distance $x_h = 30$ cm in the positive $x$-direction and a distance 10 cm in the positive $z$-direction. The components on the PCB will absorb some of the electric field that penetrates the enclosure and is incident upon the board. The PCB may, therefore, be modeled as a thin dielectric block with a reflection coefficient [18]. For this example, the reflection coefficient $\Gamma$ is uniformly distributed in the interval

$$\Gamma = [-0.91, -0.97]. \quad (43)$$

This reflection coefficient is optimized for a frequency of 1.8 GHz by changing the material parameters of the PCB. The reflection coefficient will, however, be accurate for a small frequency range around 1.8 GHz. Note that the reflection coefficient described by (43) is uncertain, and it follows a uniform distribution. The uncertainty in this input will cause there to be an uncertainty in the output.

The output $z$-component of the electric field is observed at the center of the box. An FDTD simulation was used to obtain the normalized electric field at this point using a uniform cell size of 0.01 m.

Fig. 2 shows the mean output electric field with 95% confidence intervals (CIs), as predicted by the MCM. Figures, such as this one, are extremely useful when determining the quantitative level of confidence that may be held in the results of a simulation. At 1.8 GHz, the 95% CI are $E_z = [0.418, 0.444]$ V/m. Thus, 95% of the sampled data was within about $\pm 3\%$ of the mean value.

The uncertainty in the output electric field is shown in Fig. 3. The uncertainties predicted by all three methods are in very good agreement. The uncertainty curves are compared using the FSV method over frequencies up to 3 GHz. The uncertainty predicted by the PCM is a “very good” match to the uncertainty predicted by the MCM, with a GDM of 2.3568. The MoM performs even better, and the frequency response of the uncertainty formed from the MoM and the MCM is an “excellent” match, having a GDM of 1.4755. Therefore, for this example, both efficient UA methods provide uncertainty estimates that are very close to the uncertainty formed using the MCM.

IV. EXAMPLE 2: A DIELECTRIC SPHERE

This example considers the reflection of a uniform plane wave off a dielectric sphere in free space. The incident electric field $E_y^i$ propagates in the positive $x$-direction, and is polarized in the $y$-direction with a magnitude $E_0 = 1$ V/m. The $y$-component of the backscattered field $E_y^r$ is calculated at a distance $R = 0.2$ m from the center of the sphere. This backscattered electric field is normalized relative to the input excitation to form the normalized electric field. This backscattered field may be solved analytically using the Mie series [19].

In this example, the sphere parameters are uncertain: the radius is normally distributed with a mean $\bar{a} = 0.1$ m and an uncertainty $\sigma_a = 0.005$ m, the relative permittivity is uniformly distributed in the interval $\varepsilon_r = [3.7, 4.3]$, and the relative permeability is uniformly distributed in the interval $\mu_r = [0.95, 1.05]$. These uncertain input parameters will produce an uncertainty in...
Fig. 4. Three-dimensional problem space containing a dielectric sphere. A uniform plane wave is reflected off the sphere and observed at X.

Fig. 5. Uncertainty in the normalized field backscattered from a dielectric sphere.

The normalized electric field in the frequency domain was calculated using FDTD simulations, with a uniform broad Gaussian incident plane wave and a uniform cell size of 0.005 m. Fig. 5 shows the uncertainty in the FDTD simulations calculated using the three UA methods. At the lower frequencies, the uncertainties produced by the three methods are in good agreement. However, at the higher frequencies, both the PCM and the MoM overestimate the uncertainty, when compared to the uncertainty produced by the MCM. These qualitative comparisons are confirmed using the FSV method. The PCM gives a “fair” estimate of the MCM uncertainty up to 1.02 GHz, and a “poor” estimate of the uncertainty between 1.02 and 3 GHz. The MoM does slightly better at the lower frequencies, providing a “fair” estimate of the MCM uncertainty up to 1.21 GHz. At the higher frequencies, however, the MoM performs less well, with a “very poor” estimate of the uncertainty between 1.21 and 3 GHz.

Fig. 6 shows the normalized electric field produced from an FDTD simulation using the mean input parameter values and a simulation with the sphere radius perturbed by 5 mm. The two curves in this figure have similar resonant features, but are shifted slightly in the frequency domain. At 1 GHz, the frequency response curve is less resonant in nature. Changing the radius of the sphere causes a frequency shift, which, in turn, changes the value of the normalized electric field in a quasi-linear fashion, at this frequency. Changing the radius of the sphere at a more resonant frequency (e.g., 2.71 GHz) results in a frequency shift that causes a large nonlinear change in the normalized electric field. Fig. 7 shows the relationship between the normalized electric field and the radius of the sphere at 1 and 2.71 GHz (calculated using the Mie series solution), respectively. At 1 GHz, the normalized electric field depends on the radius in a relatively linear fashion, whereas at 2.71 GHz, the normalized electric field depends on the radius in a highly nonlinear manner. Similar nonlinear relationships between the output electric field and the other uncertain inputs arise at frequencies where there is a high modal density.

In this example, the chaotic expansion used by the PCM is truncated at $P = 3$; the output is, therefore, assumed to depend linearly on the uncertain inputs. The MoM also assumes a linear relationship between the output and the uncertain inputs. At subresonant frequencies (e.g., at 1 GHz), the linear assumption is valid and the subsequent predictions of the uncertainty formed via the PCM and the MoM are similar to the uncertainty formed via the MCM. The linear assumption used by the PCM and the MoM is poorer at frequencies where the frequency response of the electric field is more nonlinear (e.g., at 2.71 GHz). This explains the poor estimations of the uncertainty produced by the PCM and the MoM at such frequencies. The resonant nature of EMC data may prevent efficient UA methods, such as the PCM and the MoM, from being used to accurately quantify the uncertainty in the frequency response of the electric field formed from CEM simulations of EMC examples. The MoM and the PCM may still be useful for quantifying the uncertainty at subresonant frequencies.
be similar despite the CEM technique used, it may be possible the uncertainty in the output of a computational scenario should computationally expensive than the MoM and the PCM. Since the nonlinear nature of EMC data. The MCM is, however, more computational simulations. The MCM, therefore, does not suffer from the uncertain inputs and the output of the computational method does not rely on any assumption on the relationship between the uncertain inputs and the output. This means that the full inner product values have to be stored at all points in the problem space. The uncertainty in the sphere radius causes the material parameter inner product values, used by the PCM, to be different at different points in the problem space. This means that the full inner product values have to be stored at each point in the problem space, requiring significantly more memory. In more complex examples, the computational memory requirements may be too large to allow the PCM to be used.

To obtain the uncertainty of these (and any other) CEM simulations, extra computational runtime is needed. This extra runtime will be significant for complex problems with many uncertain input parameters. Uncertainty budgets provide essential information to help determine whether the results of a measurement are acceptable, and should not be discounted because of the extra computational expense. This paper has investigated two efficient UA methods (the PCM and the MoM), highlighting some of the strengths and limitations of these methods.

The efficient MoM and the PCM provided poor estimates of the output uncertainty when the relationship between the output and the uncertain inputs was nonlinear. It is generally accepted that the MCM provides a benchmark method for determining output uncertainty with a good degree of accuracy. This method does not rely on any assumption on the relationship between the uncertain inputs and the output of the computational simulations. The MCM, therefore, does not suffer from the nonlinear nature of EMC data. The MCM is, however, more computationally expensive than the MoM and the PCM. Since the uncertainty in the output of a computational scenario should be similar despite the CEM technique used, it may be possible to use a computationally efficient CEM technique (such as the intermediate-level circuit model (ILCM) [20]), along with the MCM to provide an accurate estimate of the output uncertainty with relatively little computational expense. The ILCM method, for example, was able to find the shielding effectiveness of an enclosure over 3900 times faster than TLM simulations. To determine whether an accurate and efficient UA can be formed in this way, the uncertainty in the output of an EMC scenario must be shown to be relatively independent of the CEM technique used to model the scenario. This provides a promising avenue for future work in this area.

Table I displays the extra computational memory required by the PCM, compared to that required by the MCM and the MoM. In the second example, significantly, more memory is required for the PCM. The MCM and the MoM use an optimized FDTD method; the material parameter values are not stored at each point in the problem space, but a reference to the parameter value is stored. Conversely, for the PCM, material values need to be stored at all points in the problem space. The uncertainty in the sphere radius causes the material parameter inner product values, used by the PCM, to be different at different points in the problem space. This means that the full inner product values have to be stored at each point in the problem space, requiring significantly more memory. In more complex examples, the computational memory requirements may be too large to allow the PCM to be used.

V. COMPUTATIONAL REQUIREMENTS

Table I shows the computational performance of the three methods. All FDTD simulations were performed on a computer with a Pentium 4 processor running at 3.0 GHz. The table shows that the MoM requires the least amount of computational expense. The MCM requires much more computational runtime than the PCM and the MoM, highlighting the need for efficient methods of quantifying the uncertainty in CEM simulations.

<table>
<thead>
<tr>
<th>Example</th>
<th>Method</th>
<th>No. Simulations</th>
<th>Time</th>
<th>Memory</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>MCM</td>
<td>100</td>
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<td>MoM</td>
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<td>0.36 hours</td>
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<tr>
<td>1</td>
<td>PCM</td>
<td>1</td>
<td>1.8 hours</td>
<td>62 MB</td>
</tr>
<tr>
<td>2</td>
<td>MCM</td>
<td>700</td>
<td>126 hours</td>
<td>31 MB</td>
</tr>
<tr>
<td>2</td>
<td>MoM</td>
<td>4</td>
<td>0.72 hours</td>
<td>31 MB</td>
</tr>
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<td>2</td>
<td>PCM</td>
<td>1</td>
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<td>680 MB</td>
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VI. CONCLUSION

Estimates of the uncertainty in the results of CEM simulations provide the scientific community with the quantitative level of confidence that may be held in the results. In the first example of this paper, it may be concluded that there is a 95% chance that the output value lies within 3% of its mean value. It is impossible to determine this level of confidence without performing an UA.

This paper introduced three UA methods that were used to quantify the uncertainty in FDTD simulations. The novel implementation of the PCM required a modification of the FDTD algorithm. Of the three methods, the MoM was shown to be the computationally cheapest method. The PCM was shown to be computationally faster than the MCM, but required significantly more computational memory.

The MCM has been previously used to provide reliable estimates of uncertainty. The first example in this paper highlighted that the computationally cheaper MoM and PCM can give very good estimations of the uncertainty formed via the MCM. The efficient MoM and PCM, implemented in this paper, both rely on the assumption that the output of interest depends linearly on the uncertain inputs. In the second example, it was shown that this assumption is valid at subresonant frequencies, but poorer at frequencies with a higher density of resonant modes. This reflected the uncertainty estimates formed by the PCM and the MoM, for this example, which were better at subresonant frequencies. In conclusion, the MoM and the PCM may only provide moderate estimates of the uncertainty in resonant EMC data. However, the efficient methods have also been shown to work well at subresonant frequencies, for an example, with multiple uncertain inputs.

The MCM is generally accepted as being an accurate method for quantifying output uncertainties. The method does not rely on an assumed relationship between the uncertain inputs and the output of computational simulations, and therefore does not suffer from the resonant nature of EMC data. The MCM is also, however, known to be computationally expensive. It has been suggested that a fruitful avenue for future work may be to use the MCM with a computationally efficient CEM technique (such as the ILCM [20]) to provide an accurate estimate of the output uncertainty with relatively little computational expense.

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The PCM bears no relation to chaos theory, and the MoM used in this paper should not be confused with the common method

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used in CEM. Both of these methods are statistical methods that may be used to quantify uncertainty.

REFERENCES


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