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Abstract. A severe drawback to the scalar Monte Carlo (MC) method is the difficulty of introducing diffraction when simulating light propagation. This hinders, for instance, the accurate modeling of beams focused through microscope objectives, where the diffraction patterns in the focal plane are of great importance in various applications. Here, we propose to overcome this issue by means of a direct extinction method. In the MC simulations, the photon paths' initial positions are sampled from probability distributions which are calculated with a modified angular spectrum of the plane waves technique. We restricted our study to the two-dimensional case, and investigated the feasibility of our approach for absorbing yet nonscattering materials. We simulated the focusing of collimated beams with uniform profiles through microscope objectives. Our results were compared with those yielded by independent simulations using the finite-difference time-domain method. Very good agreement was achieved between the results of both methods, not only for the power distributions around the focal region including diffraction patterns, but also for the distribution of the energy flow (Poynting vector). © 2014 Society of Photo-Optical Instrumentation Engineers (SPIE) [DOI: 10.1117/1.JBO.19.11.115003]

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Introduction 1

The generation of localized light fields through high numerical aperture (NA $\ge 0.7^{1}$) optical systems has become increasingly common in the field of biophotonics, with applications ranging from optical tweezers to microscopy and spectroscopy.²⁻⁶ Yet, such systems ask for reliable simulation tools, not only for the design of the optics, but also for a better understanding of lightmatter interactions (e.g., absorption and scattering) within the focal volume. The tight focusing of beams induces diffraction patterns in the focal volume where the polarization of the illumination beam plays an important role.^{7,8} Consequently, if the tight focusing of beams is to be modeled, the consideration of the wave nature of the light becomes paramount. This can be successfully achieved by solving numerically Maxwell equations, e.g., by carrying out finite-difference time-domain (FDTD) simulations⁹ for small volumes.

The Monte Carlo (MC) algorithm¹⁰ is commonly used for calculations of the light propagation in scattering and absorbing media, and the problem of focused beams has been assessed by various groups and is still a matter of ongoing research.¹¹⁻¹⁴ We present here an alternative MC approach that enables the handling of the diffraction within absorbing media introduced by a coherent light source. Moreover, the high sampling number required for a destructive interference of a coherent source in a vectorial MC¹⁵ is rendered obsolete by this approach. In this contribution, we investigate the feasibility of including diffraction of the incident light into MC simulations in order to model tightly focused beams and Fraunhofer diffraction patterns in arbitrarily large simulation volumes in a relatively simple manner.

This paper is organized as follows: in Sec. 2, we describe how the MC simulations are combined with the so-called angular spectrum of the plane waves technique (ASPWs) to simulate the light propagation and to introduce, at the same time, our "direct extinction method." In Sec. 3, we present the results yielded by the developed method and compare them with those obtained from FDTD simulations, both in vacuo and in absorbing media. We are interested not only in the power distribution over the sample, but also in the distribution of the electric and magnetic field components and, subsequently, in the energy flow (represented by the Poynting vector) at every point. In the last section, we discuss the obtained results and present conclusions on the feasibility of our approach.

2 Theory

We consider the tight focusing of a monochromatic beam through a microscope objective (see Fig. 1). A collimated laser beam with a uniform profile and propagating along the x-direction is focused by an aplanatic cylindrical lens.¹⁶ The focal volume lies within the simulation grid. To simplify the computational effort and the simulation of larger grids, we confine our investigation here to the two-dimensional (2-D) case, so that only the x, y-plane is considered. We illustrate our purpose with two types of polarizations for the illuminating beam: first, linearly polarized along the z-axis, and second, radially polarized (within the x, y-plane). The sample and its surrounding medium are assumed to have the same index of refraction, i.e., there is no Fresnel reflection at the boundaries.

We distinguish two cases: first, the beam propagates in vacuo, and second, the beam propagates in an absorbing, yet nonscattering material.

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Fig. 1 Geometry of the simulations, as carried out in this paper. The setup is infinite in z-direction.

2.1 Direct Extinction Method within the Monte Carlo Algorithm

When dealing with light propagation problems, the MC method constitutes a well-established numerical alternative to the analytical solutions of the radiative transfer equation. ^{10,17–19} The core idea consists of stochastically modeling the light propagation with individual photon paths akin to random walks. A single photon path is composed of successive ray segments, each of which is defined by (1) a starting point, (2) a propagation direction, (3) and a segment length. Since abundant documentation can be found on the matter,^{10,15,20} we do not repeat the successive steps needed to generate such paths; instead, we emphasize the novelty of our approach. Traditionally, the initial starting point of the very first ray segment is chosen at the position of the illuminating source, or on the surface of the sample, from where the illumination is performed. Here, we propose an alternative method where the initial starting point of the rays coincides to a "target position". This position is independently sampled from a probability distribution calculated with the ASPW method and which corresponds to the probability that the ray has not interacted (by scattering or absorption) with the sample yet. We confine this study to absorbing materials as a preliminary investigation.

In this section, we detail the calculation of the probability distribution for the incident beam and the inclusion of the absorption into the ASPW method. Further, we explain how our method can be used to calculate the local Poynting vectors, as these can help to model scattering processes at a later stage. In all of these simulations, only the 2-D case is studied to simplify the computation and to decouple the polarization of the light into two orthogonal components: parallel and perpendicular. However, the consideration of the 2-D case is beyond a mere simplification of the three-dimensional (3-D) case. "In fact, both the analyses and the final results are sufficiently different in the two cases to call for separate treatments."²¹

2.2 Angular Spectrum of Plane Waves Method in Vacuo

In order to calculate the power distribution corresponding to the probability of finding a photon propagating in the simulation space, we need to know the electric field $\mathbf{E}(\mathbf{r})$ at every point $\mathbf{r} = (x, y)$. This can be obtained with the ASPW method, which is particularly well suited for problems involving arbitrarily shaped beams where the paraxial approximation fails, as in our current proposal.^{16,22} The ASPW method is a mathematical technique that allows for the representation of complex beam profiles by a coherent superposition of plane waves: this is achieved knowing that the field distribution in the focal plane is given by the Fourier transform of the field distribution in the far field, i.e., at the lens' surface (we neglect the contribution of the evanescent waves; as in our case, the focal plane is assumed to be located sufficiently afar from the lens).

Since, in a first step, the monochromatic beam is propagating in the x-direction in a nonabsorbing and nonscattering 2-D space, each plane wave in the x, y-plane is characterized by a wave vector **k**, which is defined by the wave's incidence angle θ onto the focal plane (see Fig. 1). Thus, if we consider a linearly polarized beam such that its electric field is oscillating along the z-direction ($\mathbf{E} = \mathbf{E}_z$), the angular spectrum representation of the beam yields the expression of the field

$$\mathbf{E}_{z}(\mathbf{r}) = \frac{1}{2\pi} \int_{-\infty}^{\infty} \mathbf{E}_{z}^{0}(k_{y}) \exp[j(x - x_{\text{ph}})\sqrt{k^{2} - k_{y}^{2}} + jk_{y}y] \mathrm{d}k_{y},$$
(1)

at every point **r** in the simulation space where $k = \sqrt{k_x^2 + k_y^2}$ is the wave number in the medium and $x_{\rm ph}$ is the position of the focal plane (see Fig. 2). As for $\mathbf{E}_z^0(k_y)$, it expresses the field in the focal plane and is obtained by the transformation of a collimated beam with a uniform profile truncated by an aperture incident on the cylindrical lens. We assume the field to be zero outside the aperture (Kirchhoff approximation). Further, we follow here the reasoning of Richards, Wolf, and Born,^{16,23} where the intensity law of geometrical optics allows for the calculation of the fields after passing through the lens, which leads to the formulation²⁴ of

$$\mathbf{E}(\mathbf{r}) = \frac{k}{2\pi} \int_{-\theta_{\text{max}}}^{\theta_{\text{max}}} \hat{E} \mathbf{a}(\mathbf{r}, \theta) d\theta, \qquad (2)$$



Fig. 2 Comparison between the grid sizes of FDTD and MC.

where \hat{E} denotes the amplitude of the incident electric field. Each plane wave is scaled by its respective factor of $\mathbf{a}(\mathbf{r}, \theta)$ to give a linearly polarized beam defined as

$$\mathbf{a}(\mathbf{r},\theta) = \begin{pmatrix} 0\\0\\1 \end{pmatrix} \cdot \exp[jk((x-x_{ph})\cos\theta + y\sin\theta)]\sqrt{\cos(\theta)}.$$
(3)

Note that the Kirchhoff approximation, combined with the assumption that the focal region is located sufficiently afar from the lens, corresponds to the so-called Debye-approximation.^{21,25} In order to implement a radially polarized beam, $\mathbf{a}(\mathbf{r},\theta)$ needs to be modified:²⁶

$$\mathbf{a}_{\rm rad}(\mathbf{r},\theta) = \begin{pmatrix} \sin\theta\\ \cos\theta\\ 0 \end{pmatrix} \operatorname{sgn}(\theta) \mathbf{a}_z(\mathbf{r},\theta).$$
(4)

The power $P_v(\mathbf{r}) \propto \mathbf{E}(\mathbf{r}) \cdot \mathbf{E}^*(\mathbf{r})$ leads to a probability density function $p_v(\mathbf{r})$, which expresses the probability of finding a photon at a position \mathbf{r}

$$p_v(\mathbf{r}) = \frac{P_v(\mathbf{r})}{\int_{\mathbf{r}} P_v(\mathbf{r}) d\mathbf{r}}.$$
(5)

2.3 Modeling the Absorption

Now consider the beam propagating in an absorbing material with a refractive index $\mathbf{n} = n + j\kappa$, such that its absorption coefficient is defined by $\mu_a = 2k\kappa$.

If we denote the fields of the beam *in vacuo* by $\mathbf{E}(\mathbf{r})$, the corresponding expression of the fields in the absorbing material is $\mathbf{E}^{e}(\mathbf{r}, \mu_{a})$, where the superscript "e" refers to the extinction. The influence of the medium can be represented by inhomogeneous plane waves²³ composing the beam, based on Beer-Lambert law's exponential decay (we neglect the Fresnel reflection which occurs between the nonabsorbing and the absorbing medium). The modified ASPW representation is defined as

$$\mathbf{E}^{\mathbf{e}}(\mathbf{r},\mu_{a}) = \int_{-\theta_{\max}}^{\theta_{\max}} \hat{E}\mathbf{a}^{\mathbf{e}}(\mathbf{r},\theta,\mu_{a}) \mathrm{d}\theta, \tag{6}$$

where the integrand introduces the Beer-Lambert law as

$$\mathbf{a}^{\mathrm{e}}(\mathbf{r},\theta,\mu_{a}) = \mathbf{a}(\mathbf{r},\theta) \exp\left[-\frac{1}{2}\mu_{a}\frac{x}{\cos(\theta)}\right].$$
(7)

Analogously to the case in vacuo [Eq. (5)], the power $P_a(\mathbf{r}) \propto \mathbf{E}^{e}(\mathbf{r}, \mu_a) \cdot \mathbf{E}^{e*}(\mathbf{r}, \mu_a)$ leads to a probability density function $p_a(\mathbf{r})$, which expresses the the probability of finding a photon at a position \mathbf{r} before it gets absorbed in the sample:

$$p_a(\mathbf{r}) = \frac{P_a(\mathbf{r})}{\int_{\mathbf{r}} P_a(\mathbf{r}) d\mathbf{r}}.$$
(8)

This constitutes the essence of the "direct extinction method," which can later be extended to include scattering processes. The main advantage of this approach is that the diffraction phenomena can be recovered from the sampled probability distribution without having to implement the interference between photon paths.

For simplicity, in the following part of our manuscript, we shall use the term "beam" to refer to the beam which fulfills the assumptions described in this section and whose electric field is given by $\mathbf{E}(\mathbf{r})$ *in vacuo* and by $\mathbf{E}^e(\mathbf{r}, \mu_a)$ in an absorbing medium.

2.4 Ray Direction: Poynting Vector

As already mentioned, the present study is confined to nonscattering materials, i.e., the MC simulations do not need to propagate the photon paths, and the sampling from the aforesaid probability distribution suffices. However, if scattering processes are to be modeled, another key information besides the "first interaction point" needs to be known: the first path segment's propagation direction.

To this end, we may utilize the energy flow or the energy flux density given by the local time-averaged Poynting vector $\boldsymbol{S}(\boldsymbol{r})$

$$\mathbf{S}(\mathbf{r}) \propto \operatorname{real}\{\mathbf{E}^{e}(\mathbf{r}) \times \mathbf{H}^{e*}(\mathbf{r})\},$$
(9)

where the magnetic field \mathbf{H}^{e} at a point **r** is calculated similarly to the electric field with the modified ASPW method to give

$$\mathbf{H}^{\mathbf{e}}(\mathbf{r},\mu_{a}) = \int_{-\theta_{\max}}^{\theta_{\max}} \mathbf{b}^{\mathbf{e}}(\mathbf{r},\theta,\mu_{a}) \mathrm{d}\theta, \qquad (10)$$

where \mathbf{b} denotes the contribution of the single plane waves. For the linearly polarized beam we obtain

$$\mathbf{b}^{\mathbf{e}}(\mathbf{r},\theta) = \hat{H} \begin{pmatrix} \sin \theta \\ \cos \theta \\ 0 \end{pmatrix} \mathbf{a}_{z}^{\mathbf{e}}(\mathbf{r},\theta).$$
(11)

 \hat{H} denotes the amplitude of the incident magnetic field. As for the radially polarized beam we obtain

$$\mathbf{b}_{\rm rad}^{\rm e}(\mathbf{r},\theta) = \hat{H}\,{\rm sgn}(\theta)\mathbf{a}^{\rm e}(\mathbf{r},\theta). \tag{12}$$

Note that for $\mu_a = 0$, the magnetic and electric fields \mathbf{E}^e and \mathbf{H}^e are reduced to the fields \mathbf{E} and \mathbf{H} *in vacuo*, respectively.

3 Simulations and Results

3.1 *Principle: in Vacuo*

We simulated the 2-D version of the situation sketched in Fig. 1: the beam (as described in the previous section) propagating in the *x*-direction is focused by an aplanatic microscope objective, so that the divergence angle is $\theta_{max} = 65$ deg, which lies beyond the paraxial limit. This corresponds to a numerical aperture of NA = 0.9 in air or NA = 1.2 in water, and is relatively common for modern confocal microscope objectives. As an example, we consider the monochromatic beam with a wavelength of $\lambda = 1 \ \mu$ m. Note that by scaling, these results are also valid for any other wavelength.

The MC simulations were carried out in a square-shaped sample, with 200 bins in the x- and y-directions, with bin sizes of $\Delta x = \Delta y = \lambda/25$. As in the ASPW model, Fresnel reflection at the sample interfaces was neglected, i.e., the sample's refractive index was assumed to be equal to that of the surrounding medium ($n_{\text{sample}} = n_{\text{air}} = 1$). Since we are dealing with nonscattering materials, the MC simulations are confined to sampling "target positions" from the normalized probability distribution $p_v(\mathbf{r})$ calculated with the ASPW method, where 41 plane waves were superimposed. In vacuo, $p_v(\mathbf{r})$ simply expresses the probability of finding a photon at a position **r**. The sampling is achieved with the rejection method. First, a pseudorandom position \mathbf{r}_i is drawn. (Pseudorandom numbers were generated with the Mersenne Twister²⁷ algorithm.) The corresponding probability density $p_v(\mathbf{r}_i)$ at that position is compared with a pseudorandom number $\xi \in (0,1]$: If the number is such that $\xi > p_v(\mathbf{r}_i)$, the position is rejected and the sampling is repeated.

In order to assess the accuracy of our results, we conducted independent FDTD simulations (more details on these simulations can be found elsewhere^{9,24,28}). These simulations numerically solve the discretized time-dependent Maxwell's equations of electrodynamics to propagate the electric and magnetic fields within a grid using a leap frog algorithm. As input, we need to specify for each plane wave its (1) incidence angle θ on the sample, together with its subsequent starting position $\mathbf{r}_0 = (x = 0, y)$, and (2) its electric field at \mathbf{r}_0 . To remain consistent with the MC simulations, we propagate 41 plane waves with incidence angles lying in the range $[-\theta_{max}, +\theta_{max}]$. For a given incidence θ , an expression of the electric field at \mathbf{r}_0 is obtained from the ASPW calculations [i.e., from Eq. (2)]. All FDTD simulations throughout the work were performed with a resolution of $\lambda/25$.

Results obtained with MC and FDTD simulations for a linearly polarized beam propagating *in vacuo* are shown in Fig. 3. In this paper, all simulation results are displayed after normalization by the probability density value at the focal point $(x = x_{ph}, y = 0)$ of the simulation *in vacuo*, denoted by $p_{v,0}$ and $p'_{v,0}$ for the linearly and radially polarized beams, respectively. Both types of simulations were performed with the same grid dimensions. The agreement between the two sets of data serves as a verification to the implementation of both the MC and the FDTD algorithms.

Besides, as the FDTD method accounts for the nearfield interactions, we can verify here the formation of the Fraunhofer diffraction patterns in the vicinity of the focal plane. However, we stress here the main constraint inherent to the MC technique: the number of samplings should be large enough to provide reasonable convergence.

3.2 In Absorbing Media

As a following step, we compare the MC and FDTD simulations in the presence of absorption. This would be of interest, for instance, to investigate the excitation of fluorophores with confocal fluorescence systems. Again, the MC simulations were performed with a square-shaped sample of dimensions $8\lambda \times 8\lambda$, such that the focal plane of the focused beam is located at $x_{ph} = 4\lambda$.

Four types of samples were considered, each sample having a different absorption coefficient $(16\lambda)^{-1} \le \mu_a \le (2\lambda)^{-1}$. Note that despite the large absorption values, we omit here re-emission, heat transfer, denaturation, etc., since these do not contribute to our present study. As opposed to the choice of these high absorption coefficients, it is also possible to simulate smaller values, but that requires a much larger grid in order to show an effect on the focus. However, an increased grid size is disadvantageous for the performance of the FDTD simulation.

We proceed according to our extinction method, by sampling "target positions" from $p_a(\mathbf{r})$ instead of $p_v(\mathbf{r})$. These "target positions" bear a slightly different significance when absorption is involved: $p_a(\mathbf{r})$ represents the probability to find a photon in the sample before it gets absorbed along its path.



Fig. 3 Comparison of the probability distributions yielded by the Monte Carlo (MC) simulations (the number of sampled "target positions" are 10^6 and 10^9 on (a) and (b), respectively) and the finite-difference time-domain (FDTD) simulations for the linearly polarized beam focused *in vacuo* in the *x*-direction. These distributions are displayed here in logarithmic scale over the range $8\lambda \times 8\lambda$. The MC simulations are shown for $y/\lambda > 0$ and FDTD simulations for $y/\lambda \le 0$ for both figures. The curves are cross sections of the probability profiles at the focal plane $x = x_{ph} = 4\lambda$.



Fig. 4 Three different FDTD grid sizes displayed in logarithmic scale over the range $8\lambda \times 8\lambda$ for the linearly polarized beam, from (a) $x = 8\lambda$ and $y = 8\lambda$, (b) $8\lambda \times 16\lambda$, and (c) $8\lambda \times 32\lambda$.

In the FDTD simulations, the absorption in the sample is taken into account by increasing the imaginary part of the refractive index: $(64\pi)^{-1} \le \kappa \le (8\pi)^{-1}$. The grid spacing is preserved, but the overall dimensions of the simulation grid need to be modified. This is shown in Fig. 2. Precaution is needed when comparing the MC simulation results (based on a propagation space infinite in the y-direction) with those yielded by FDTD simulations: in the case of larger divergence angles θ_{max} , the inclusion of absorption in the medium where the beam propagates requires an adjustment of the finite FDTD simulation grid. In the figure, the darker region corresponds to the MC simulation area. The symbol x_{ph} indicates the position of the focal plane. The FDTD grid needs to be wide enough so that all plane waves start getting attenuated at x = 0; this principally concerns plane waves incident on the focal plane with a highly oblique angle θ . The dashed area corresponds to the extra computational space: the choice of the additional space naturally affects the outcome of the simulations. The interaction between the incident plane waves (required by ASPW) and the absorbing medium requires an infinite boundary. However, FDTD can only calculate finite volumes.

An optimal grid size can be evaluated as shown in Fig. 4. Here, a larger grid space in the FDTD simulations compared with the grid considered in the MC simulations is evaluated. We show here the power distribution obtained with three different FDTD grid sizes for the linearly polarized beam focused in an absorbing material with $\mu_a = (2\lambda)^{-1}$ in the *x*-direction. The grid size in the FDTD simulations needs to be large enough to provide correct results by accounting for the absorption of plane waves with a wide angular incidence on the focal plane. The figure in the center and on the right reveals that the grid size needs to be above a certain threshold for the results to converge.²⁹

Simulation results obtained with the four types of samples are shown in Fig. 5 and more detailed profiles from the same results are shown in Fig. 6. It can be seen that good agreement is achieved between the MC and the FDTD simulations. Nonetheless, closer inspection reveals the appearance of discrepancies, both close to the sample's illumination face and in the focal region (the latter case is not shown here but was observed, for instance, in the case of beams focused with a divergence angle of $\theta_{\text{max}} = 25$ deg). Since similar discrepancies have been observed no matter what divergence angle $\theta_{\rm max}$ was taken for the beam, the most plausible explanation for the differences is the presence of Fresnel reflection in the FDTD simulations. In the FDTD model, the plane waves propagate from air into a sample, whose refractive index has a nonzero imaginary component. However, as mentioned earlier, the MC simulations were implemented assuming that there is no refractive index mismatch along the plane waves' path.

3.3 Direction of the Poynting Vectors

In the last step of our investigation, we compared the orientation of the Poynting vectors $S(\mathbf{r})$ as calculated with both MC and



Fig. 5 Comparison of the probability distributions displayed here in logarithmic scale over the range $8\lambda \times 8\lambda$ and yielded by the MC simulations (with 10⁹ "target positions" sampled) and the FDTD simulations (grid size $8\lambda \times 32\lambda$) for the linearly polarized beam focused in the *x*-direction in increasingly absorbing media. From (a) to (d): $\mu_a = (16\lambda)^{-1}$, $(8\lambda)^{-1}$, $(4\lambda)^{-1}$ and $(2\lambda)^{-1}$ (the corresponding values for the imaginary component κ of the material's refractive index are $(64\pi)^{-1}$, $(32\pi)^{-1}$, $(16\pi)^{-1}$, and $(8\pi)^{-1}$, respectively). The curves are cross sections of the probability profiles at the focal plane $x = x_{ph}$ (MC in blue and FDTD in green).



Fig. 6 Quantitative representation of the results shown in Fig. 5. The probability profiles corresponding to the different absorbing materials are cross sections close to the illumination face (a), at $x = 0.01\lambda$ and in the focal plane at $x = x_{ph} = 4\lambda$ (b).

FDTD simulations. In the MC method, the Poynting vector is calculated according to Sec. 2.4, averaged over a bin of subwavelength size $\Delta x \times \Delta y$. On the other hand, in the FDTD simulations, the Poynting vector is determined at a certain grid point (x, y) from the resulting electric and magnetic fields at the end of the simulation, see Eq. (9).

The results obtained by focusing a linearly polarized beam are shown in Fig. 7. Again, very good quantitative agreement is achieved between the two datasets. Yet, differences can be seen in the vicinity of the focal plane, in particular in the darker regions of the probability distribution where destructive interferences occur. In those regions, the orientation of the Poynting vector might vary considerably over a bin, in which case the averaging with the MC method does not yield identical results to the FDTD simulations. In other regions where discrepancies are not visible, the orientation of the Poynting vector remains relatively constant over a bin, and thus the averaging produces identical results to the FDTD simulations. Moreover, the effect of the aforementioned Fresnel reflection included in the FDTD calculations and neglected in the MC simulations should not be excluded in the interpretation of the differences. Avoiding the mismatch of the imaginary part of the refractive index in FDTD by adjusting the refractive index in the outer regions of the grid is a nontrivial task and would require further development of the FDTD algorithm, which is beyond the scope of the present feasibility study.

Additionally, results yielded by the focusing of a radially polarized beam are shown in Fig. 8: The corresponding Fraunhofer diffraction pattern noticeably differs from that of the previous example. As for the Poynting vectors in Fig. 8, the remarks on Fig. 7 still hold true. Here the observed spacial distribution depends on the resulting electric field from both E_x and E_y . Furthermore, the tight focusing of the radially polarized beam leads to a strong longitudinal electric field component in



Fig. 7 Direction of the normalized Poynting vectors **S**(**r**) plotted over probability distributions displayed here in logarithmic scale over the range $3\lambda \times 3\lambda$, and calculated with both MC simulations (10^9 "target positions" sampled) and FDTD simulations (grid size of $8\lambda \times 32\lambda$). (a) The linearly polarized beam is focused *in vacuo*. (b) The same beam is focused in an absorbing medium, where $\mu_a = (2\lambda)^{-1}$. In MC simulations, the binning size was kept at $\Delta x = \Delta y = \lambda/25$.



Fig. 8 Radially polarized beam is focused *in vacuo*.(a) Distribution of the non-normalized Poynting vectors **S**(**r**) plotted over probability distribution displayed here in logarithmic scale over the range $4\lambda \times 4\lambda$, and calculated with FDTD simulations (grid size of $8\lambda \times 32\lambda$). (b) Same distribution of the normalized Poynting vectors obtained with both MC simulations (10^8 "target positions" sampled) and the FDTD simulations, plotted over probability distribution displayed here in logarithmic scale over the range $8\lambda \times 8\lambda$.

the *x*-direction in the focal region, and the Poynting vector vanishes [as can be seen from the FDTD simulations of Fig. 8(a)].^{30–32}

4 Discussion and Conclusions

In this contribution, we have investigated the feasibility of modeling the tight focusing of beams with MC simulations based on the "direct extinction method," where "target positions" for the photon paths are sampled from probability distributions separately calculated with the ASPW technique. The simulations present an efficient tool to examine some of the wave properties of the light (Fraunhofer diffraction pattern, polarization state, and the direction of the energy flow) in the focal region of a high-aperture system, and unlike the FDTD simulations, the MC method presented here does not impose limitations on the dimensions of the observed sample.

Naturally, the accuracy of the results depends not only on the spatial resolution, but also on the number of sampled "target positions" and the number of plane waves used. While the FDTD grid has to be fine enough to avoid numerical artifacts, the MC binning can be broadened to a limit given by the theorem of Nyquist-Shannon ($<\lambda/2$). Although we have chosen here to model the focusing of a collimated beam with a uniform profile in the angular domain, more intricate beam profiles (Gaussian, Bessel, etc.,) can be easily implemented.²⁹

Likewise, we can conclude the following from our proof of concept *in vacuo* and in absorbing media: (1) the "direct extinction method" using inhomogeneous plane waves yields a very good approximation when treating problems involving absorption; (2) the assumptions adopted in the FDTD simulations (as explained in Sec. 3.2) work equally reasonably well as shown by the very good agreement with the MC results despite the numerical challenges involved. Certainly, including Fresnel reflection phenomena and extending our model to the 3-D case would be of particular importance in a next step.

Another significant improvement would be the treatment of scattering processes within the medium in which the beam is focused. A preliminary investigation has been undertaken here by calculating the direction of the energy flow in the vicinity of the focal region. This highlights a meaningful advantage of our method: We have shown that our alternative MC approach can also account for waves that show longitudinal components (nonTEM, as for example in Fig. 8). The modeling of such longitudinal components is part of the limitations of the existing vectorial MC programs: this necessitates the development of new adequate models, as initiated by Azzam.³³

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