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# Rigorous analysis of numerical methods: a comparative study

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**Abstract** For any photonic device simulation, the accuracy of the numerical solution not only depends on the methods being used but also on the discretization parameters used in that numerical method. In this work, Finite Element Method and Finite Difference Time Domain Method based on Maxwell's equations were used to simulate optical waveguides and directional couplers. As the solution accuracy may also depend on the index contrast used in such photonic devices, the characteristics of low-index contrast Germanium doped Silica and high-index contrast Silicon Nanowire Waveguides were analyzed, evaluated and benchmarked. Numerical results to benchmark Directional Couplers are also reported in this paper.

**Keywords** Waveguides, couplers, and arrays · Finite element methods · Integrated optics

## 1 Introduction

The propagation of light through optical guided wave devices can be characterised by using various modelling techniques. These techniques can be classified into analytical methods and numerical methods. Analytical methods, if possible, a solution may be

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obtained by solving the basic electromagnetic equations. In contrast, numerical methods are computational schemes or models that can be applied to a spectrum of problems by modifying the basic model to fit the problem. Be it an analytical or numerical, an accurate modelling of optical waveguides and devices is important.

However, due to the complex nature for modern optical devices, the use of analytical methods is restricted to only simple structures. Even for two-dimensional optical waveguide structures, analytical solutions are not possible and some approximations have to be made (Rahman and Agrawal 2013). Instead, existing and improved numerical methods and techniques are receiving wider attention for modeling optical components and devices, such as the Method of Moments (MoM) (Garcia et al. 2002; Hagness et al. 1997), the Finite Element Method (FEM), Beam Propagation Method (BPM), Finite Difference Method (FDM), the Finite Difference Time Domain Method (FDTD) (Luebbers 1994), the Transmission Line Modeling method (TLM), and the Time Domain Integral Equation (TDIE) techniques (de Electromagnetismo et al. 1992).

## 2 Theory

### 2.1 Wave equation

Modal analysis of optical waveguides implies the process of finding the propagation constants and the field profiles of all the modes that a waveguide can support. To obtain these propagation characteristics, solutions of the well-known Maxwell's equations given below, are necessary along with the satisfaction of the associated boundary conditions (Rahman and Agrawal 2013):

$$\nabla \times \mathbf{E} + \frac{\partial \mathbf{B}}{\partial t} = 0 \quad (1)$$

$$\nabla \times \mathbf{H} - \frac{\partial \mathbf{D}}{\partial t} = \mathbf{J} \quad (2)$$

In an isotropic lossless medium with no source ( $\mathbf{J} = 0, \rho = 0$ ), with constant permeability  $\mu = \mu_0$ , by eliminating the magnetic flux density in and the electric flux density components for Maxwell's Eqs. (1) and (2) can be written as:

$$\nabla^2 \mathbf{E} + k^2 \mathbf{E} = 0 \quad (3)$$

$$\nabla^2 \mathbf{H} + k^2 \mathbf{H} = 0 \quad (4)$$

where, the wavenumber,  $k$  (rad/m) is given as;

$$k = \omega \sqrt{\epsilon \cdot \mu_0} \quad (5)$$

Equations (3) and (4) are called Helmholtz wave equations (März 1995) for homogeneous media.

### 2.2 Analytical method

Analytical solution is only possible for planar (slab) optical waveguides by solving a resultant transcendental equation. For an optical waveguide with two-dimensional

confinement, analytical solution is not possible. Marcatili's Method was one of the first semi-analytical approximation methods developed for the analysis of buried waveguides and couplers (Marcatili 1969). This method works well in the regions far from cut-off but does not provide a satisfactory solution close to cut-off region (Chiang 1994).

Subsequently, the Effective Index Method (EIM) was proposed by Knox and Toullos in 1970 (Knox and Toullos 1970) as an extension to the Marcatili's method that became one of the most popular methods in the 1970s for the analysis of optical waveguides whereby the rectangular structure is replaced by an equivalent slab with an effective refractive index obtained from another slab. The disadvantage of this method is that it does not give good results when the structure operates near cut-off region. However, the simplicity and speed of the method have encouraged many researchers to search for different approaches to improve the accuracy of the EIM, which subsequently led to many different variants of the EIM to be developed including the EIM based on linear combinations of solutions (Chiang 1986) or the EIM with perturbation correction (Chiang et al. 1996).

## 2.3 Numerical methods

On the other hand, the Finite Difference Method, Finite Element Method, Beam Propagation Method, and the Finite Difference Time Domain Method are the popular numerical analysis methods used in many Engineering simulations. With continuous improvement of computational power at a reduced price, made these numerical methods more versatile, accurate and cost-effective.

### 2.3.1 Finite difference method

The Finite Difference Method (FDM) is one of the oldest and perhaps the most commonly used numerical techniques, until recently, in analysing optical waveguide problems. Its application to the modelling of optical waveguides dates from the early eighties, originally evolving from previous finite difference models for metal waveguides (Davies and Muilwyk 1966). The finite difference method discretizes the cross section of the waveguide that is being analysed and it is therefore suitable for modelling inhomogeneous media and complicated boundaries. In the FDM, it is necessary to define the finite cross section by enclosing the dielectric guide in a rectangular box, with the side walls as either electric or magnetic walls and the field at these boundaries are assumed to be very small. However, if leakage losses need to be calculated, these hard boundaries can be replaced by Perfectly Matched Layer (PML) (Berenger 1994). The enclosed cross section is divided into a rectangular grids allowing for the material discontinuities only along mesh lines (Bierwirth et al. 1986). The nodes are placed on mesh points so that each node can be associated to a maximum of four or eight neighbouring nodes and each node can be of one or more field variables depending on vector, semi-vectorial or scalar wave equations that can be approximated in terms of the fields at the neighbouring nodes of the mesh. Taking into account the boundary conditions of the electric and magnetic components at the dielectric interfaces, an eigenvalue problem is generated which can be solved in order to obtain the modal propagation constants and their modal field profiles. The accuracy of the method depends on the mesh size, the assumed nature of the electromagnetic field (scalar, semi-vectorial or vector) and the order of the finite difference scheme used. Often uniform mesh is used which can result in a very large number of nodes and large matrices and therefore may be computationally less efficient, particularly when dimensions of the sub-regions can be widely different, such as a graphene clad silicon nanowire.

### 2.3.2 Finite element method

The Finite Element Method (FEM) is a numerical technique for solving a wide variety of engineering problems, including computational electromagnetics.

In FEM formulation, the domain of interest is divided into many discrete elements, often triangular, which is also the principal characteristic of the method. Each element is then possible to have different material properties in terms of its relative permeability and relative permittivity, thus making the dielectric material lossy, anisotropic, or non-linear, if necessary.

The accuracy of the method also depends upon the mesh, although a finer mesh across the whole domain may yield accurate results but at the cost of increased computing time. A finer mesh can be used in areas, where the field will have a rapid variation and/or higher magnitudes and a much coarser mesh in those areas where there is little variation or negligible field. Elements should not contain physical boundaries, i.e., there should be no abrupt change in property (e.g., refractive index) within the confines of an element even though the property may change from element to element. Symmetrical domains should have symmetrical meshes as well, although whenever possible it may be better to take advantage of the symmetry in the waveguide by using the appropriate boundary conditions along the line of symmetry. The variational procedure is applied to this functional by the way of the stationary requirement, from which the Euler equation is derived and which corresponds to the wave equation. We have used a magnetic-field based formulation as shown in Eq. (6), which is then integrated over the domain (Rahman and Davies 1984).

$$\omega^2 = \frac{\int (\nabla \times \mathbf{H})^* \cdot \hat{\epsilon}^{-1} \cdot (\nabla \times \mathbf{H}) d\Omega}{\int \mathbf{H}^* \cdot \hat{\mu} \cdot \mathbf{H} d\Omega} \quad (6)$$

where,  $\omega$  is the angular frequency of each waveguide mode,  $\Omega$  is the waveguide cross-section and,  $\hat{\epsilon}$  and  $\hat{\mu}$  are the permittivity and permeability tensors of the loss-free material, respectively.

The Rayleigh-Ritz approach can be applied to this (variational) formulation by using an interpolation function of the elements used to discretise the domain. The polynomial function, which approximates the field, should remain unchanged under a linear (or higher order) transformation from one co-ordinate system to the other. All the element contributions should then be combined to form a global matrix. The resultant system of (matrix) equations or rather matrix eigenvalue equation given in Eq. (7) (Rahman and Davies 1984), should be solved using an appropriate matrix solver, as the matrices are generally very sparse.

$$[A]\{\mathbf{x}\} - \lambda[B]\{\mathbf{x}\} = 0 \quad (7)$$

where,  $[A]$  is a complex Hermitian matrix and  $[B]$  is a real symmetric and positive-definite matrix,  $\lambda$  is the eigenvalue and this can be taken as  $k_0^2$  or  $\beta^2$  depending on the formulation used. However, for a loss-less optical waveguide, by considering  $H_z$  90° out of phase with the transverse components, the  $[A]$  and  $[B]$  matrices can be transformed to Real and Symmetric and its solution would then be easier.

### 2.3.3 Beam propagation method

The Beam Propagation Method (BPM) describes the evolution of the total field propagating along a guided-wave structure and it is the most widely used tool in the study of

light propagation in longitudinally varying waveguides such as tapers, Y-junctions, and bends. The beam propagation method was first applied to optoelectronics in 1980 (Feit and Fleck 1980) and the solutions for the optical waveguides can be made to generate mode-related properties such as propagation constants, relative mode powers and group delays with high precision and considerable flexibility. The first reported BPM was based on the Fast Fourier Transform (FFT) and only solved the scalar wave equations under paraxial approximation. Therefore the FFT-BPM is only suitable for the case of weakly guiding structures, neglecting the vectorial properties of the field. Several numerical algorithms to treat the vectorial wave propagation (vector BPM) using the finite difference method, have been reported (Chung et al. 1991; Huang et al. 1992a; Huang and Xu 1992b). The VBPMs are capable of simulating polarized or even hybrid wave propagation in strongly guiding structures. Subsequently, the finite element method has been utilised to develop BPM approaches. A unified finite element beam propagation method has been reported (Tsuji and Koshiba 1996) for both TE and TM waves propagating in strongly guiding longitudinally varying optical waveguides. Obayya et al. (2000) has reported a full-vectorial BPM algorithm based on the finite element method to characterise 3-D optical guided wave devices.

Although imaginary distance BPM (Obayya et al. 2000) can find modes in a uniform optical waveguide, but being a 3-dimensional method, these are numerically costlier than 2-D FDM or FEM based modal analysis approaches.

### 2.3.4 Finite difference time domain method

However, if the optical structure contain strong discontinuity or nonlinearity or time dependent excitation, a time-domain approach would be necessary.

Kane Yee published the first paper in May 1966 to describe Maxwell's equations in the equivalent set of finite difference equations, thus showing the discretisation of space and time (Yee 1966). This led to further investigation of time-varying media and also formed the basis of the Finite Difference Time Domain (FDTD) method. Yee's approach applies a simple, second order accurate central-difference approximations for the space and time derivatives of the electric and magnetic fields directly to the respective differential operators of the curl equations. The algorithm proposed by Yee solves for both electric and magnetic fields using the coupled Maxwell's equations instead of solving individual components (electric or magnetic) alone with a wave equation, which is similar to the combined-field integral equation of the Method of Moments. Thus, creating a more robust solution in a straight forward manner in which the electric and magnetic properties of the material can be modeled. This method divides the three-dimensional geometry into cells to form a grid, called Yee cell.

Yee's leapfrog algorithm solves for all the  $\mathbf{E}$  components at a given time-step based on previously computed and stored in memory  $\mathbf{H}$  data. Then, the magnetic field is computed at the next time-step using the  $\mathbf{E}$  data just computed. The process is repeated until the time-stepping is concluded. As the algorithm runs, at each time-step the system of the electric and magnetic field components are updated based on the system of equations, which is fully explicit. In this case, there is no need to solve a system of linear equations, and the required computer memory and time is proportional to the size of the computational domain, which is based on the size of the photonic structure modelled. The computational domain is the space under consideration, where the simulation will be performed. Since the FDTD method is an explicit scheme, there is a limit on the time step  $\Delta t$  to ensure stability in the algorithm. The choice of  $S$ , called numerical stability or the Courant–Friedrichs–

Lewy factor, is then essential to minimise numerical instability, numerical dispersion and the numerical phase-velocity discontinuity. This stability factor for three dimensional geometry is then given by Taflove and Hagness (2005),

$$S = c \cdot \Delta t \sqrt{\frac{1}{(\Delta x)^2} + \frac{1}{(\Delta y)^2} + \frac{1}{(\Delta z)^2}} \quad (8)$$

where,  $c$  is the wave propagation speed,  $\Delta t$  is the time step,  $\Delta x$ ,  $\Delta y$  and  $\Delta z$  are the space increments in the x, y and z directions, respectively, and the stability condition as,  $S < 1$ .

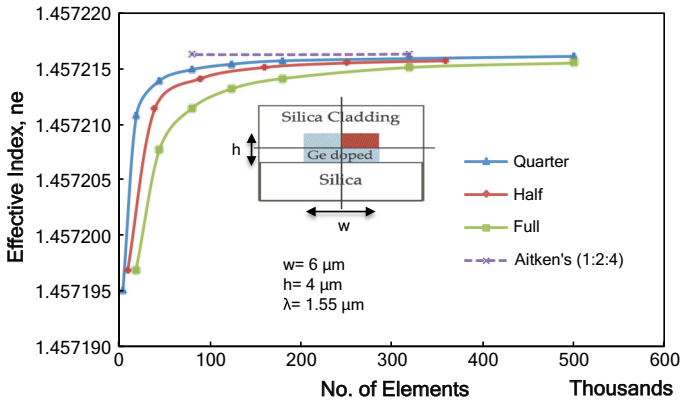
Although FDTD is a very powerful numerical approach but being a 4-dimensional approach (with added time), it is numerically much more expensive than the methods described earlier.

### 3 Numerical results

#### 3.1 Benchmarking

To consider a low-index contrast structure, first a waveguide with Germanium doped Silica core is studied, where cladding is pure Silica with refractive index of 1.44427. The core width is set at 6  $\mu\text{m}$  and height at 4  $\mu\text{m}$ . The index difference of the core with the cladding is considered to be 1.5 %, which enables us to investigate the behaviour of a low index contrast material. Using the **H**-field FEM (Rahman and Davies 1984), the numerical analysis has been carried out to find the effective index of the fundamental mode with different mesh divisions. The waveguide structure is simulated as a full waveguide and also exploiting onefold symmetry and twofold symmetry. The simulations are carried out at 1550 nm operating wavelength.

Figure 1 shows the variation of effective index with the number of elements being used for the  $H_{11}^y$  mode. Here, effective index is defined as the ratio of propagation constant,  $\beta$ , to the wavenumber,  $k$ . It can be clearly observed that the effective index increases with the increase in number of elements to a certain level and then settles asymptotically. Initially,



**Fig. 1** Variation of effective index with the number of elements for the fundamental quasi-TE ( $H_{11}^y$ ) mode of the Ge doped silica waveguide



simulation was conducted for full waveguide structure and the result of the effective index is shown by a green line. As often there may be a limitations on the number of elements that can be used or computational resources needed, it should be noted that, if a structure has onefold or twofold symmetry, then this symmetry can be exploited. In this case, to show the advantages of exploiting symmetry, results for both the cases are also shown in Fig. 1. As such, the effective index obtained for half symmetry is higher than that of full structure, shown by a red line. This shows that for a given number of elements, results using onefold symmetry is more accurate than that of considering the full structure. Similarly, in case of a twofold symmetry, it is even higher, shown by a blue line in this figure, compared to the other two lines. This clearly proves that whenever symmetry condition(s) exists, this can be exploited to improve the solution accuracy.

It should also be noted, sometimes several modes can have the same or very close eigenvalues. In that case, these modes can degenerate and eigenvectors can be mixed up. If symmetry condition is available, and used, mode degeneration can be avoided.

It is known that various extrapolation techniques can also be used to improve the solution accuracy further. Amongst them, Aitken's extrapolation is a powerful one but this requires 3 successive solutions using a fixed geometric ratio of the mesh refinements, as given below (Rahman and Davies 1985).

$$x_{\infty} = x_{r+1} - \frac{(x_{r+1} - x_r)^2}{x_{r+1} - 2x_r + x_{r-1}} \quad (9)$$

where,  $x_{r-1}$ ,  $x_r$ ,  $x_{r+1}$  are the results for three successive mesh refinements, and  $x_{\infty}$  is the extrapolated result.

To test the convergence, the results obtained from twofold symmetry is used in Eq. (9), considering the mesh division ratio of 1:2:4. The result obtained is shown by the purple dotted line in Fig. 1. In order to benchmark our results, the commercially available packages, COMSOL, FIMMPROP and the FDTD based Lumerical have been used for the simulation of the same waveguide structure.

Table 1 shows the effective index of Ge doped Silica core for fundamental quasi-TE,  $H_{11}^y$  mode. The effective indices obtained using COMSOL, FIMMPROP, Lumerical and the Aitken-extrapolation method are also tabulated in Table 1.

The result shows effective indices are accurate upto the 5th decimal, for most of the approaches used here. The normalised propagation constant, 'b' variation with mesh size for Germanium (Ge) doped Silica is much less with 0.001 only, where 'b' is given by Marcatili (1969),

$$\mathbf{b} = \frac{(n_e)^2 - (n_s)^2}{(n_g)^2 - (n_s)^2} \quad (10)$$

where,  $n_e$  is the effective index of the waveguide, and  $n_s$  and  $n_g$  are the refractive indices of Silica cladding and the doped-Silica core, respectively.

**Table 1** Effective index of Ge doped silica using various numerical methods

FEM (quarter structure)	Aitken's	COMSOL	FIMMPROP	Lumerical
1.4572161	1.4572163	1.457214	1.4572164	1.457237

Most of the optical waveguides are open-type structure and not confined inside a box. However, all the numerical method considers a finite computational region for analyses and at the computational boundary an artificial boundary condition is introduced depending on the natural boundary condition of the formulation used. These boundary conditions can be classified as (Itoh 1989),

$$\phi = 0 \text{ Homogenous Dirichlet} \quad (11)$$

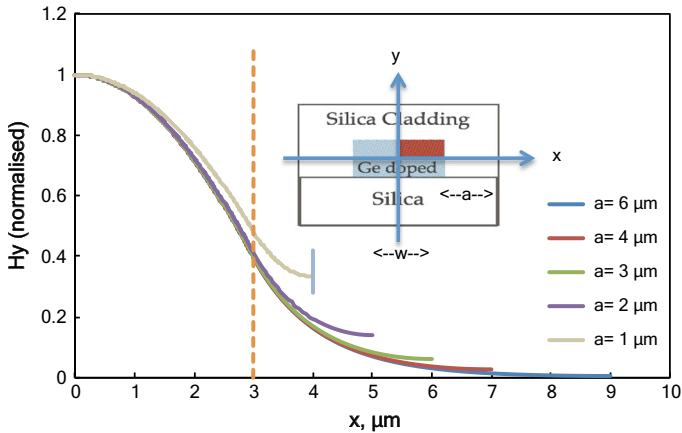
$$\phi = \mathbf{k} \text{ Inhomogeneous Dirichlet} \quad (12)$$

$$\frac{\partial \phi}{\partial t} = 0 \text{ Homogenous Neumann} \quad (13)$$

where,  $\phi$  can be electric or magnetic field and  $\mathbf{k}$  is a prescribed constant value. The vector  $\mathbf{H}$ -field formulation described in Eq. (6) has the natural boundary condition of an electric wall, i.e.  $\mathbf{n} \cdot \mathbf{H} = 0$ , where,  $\mathbf{n}$  is the unit vector normal to the surface. Therefore, when this formulation is used, there is no need to force any boundary condition on conducting guide walls.

Earlier a powerful approach, infinite element was introduced to extend the domain of the field representation (Rahman and Davies 1984). So, it is important to study the effect of using infinite element in which the computational domain can be extended in both the transverse directions.

Figure 2 shows the  $H_y$  field variation along the x-direction without infinite element. The  $H_y$  field variation obtained so far for the different waveguide width is converted to a normalised form to compare them easily. Here, 'a' is shown as the distance of the right hand side boundary from the edge of the guide. When 'a' is reduced, Electrical wall comes closer to the waveguide core and influences the modal field. For  $H_{mn}^y$  mode, the vertical metal (electric) wall imposes Neumann boundary condition on the dominant  $H_y$  field. The field plots in Fig. 2 show the case of Neumann boundary condition with  $\partial H_y / \partial x = 0$  at the computational wall. The introduction of metal wall on the side is equivalent to a multicore periodic waveguide array. In this case, (for  $H_{11}^y$  mode), as the electric wall introduces even-



**Fig. 2**  $H_y$  field variation along width of waveguide for Ge doped Silica with orthodox computational boundary

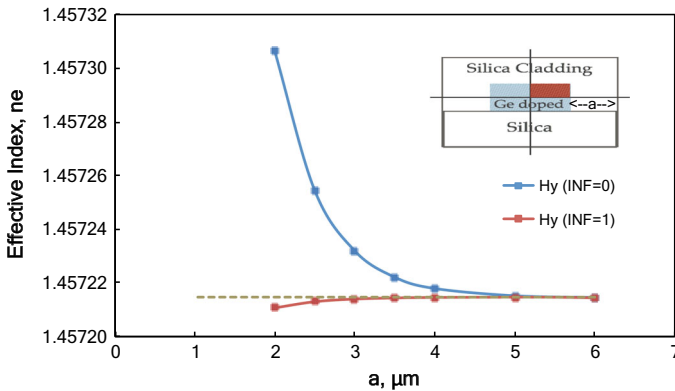
like supermode field profile, so introduces error in mode profile and this also cause the effective index to increase as shown in Fig. 3.

Figure 3 shows variation of effective index with respect to the distance,  $a$ , the distance between the waveguide core edge to the boundary wall of the computational domain, with and without infinite element. The blue line in the graph represents the effective index variation without infinite element, whereas, red line with the infinite element. In this case, effective index variation is more stable with the infinite element, even though the boundary wall is brought much closer to the core, compared to that of without infinite element as the percentage change in such variation is very less. It is being observed that  $\Delta n$  without infinite element is 0.0062856 %, whereas, with infinite element is 0.0002745 % only.

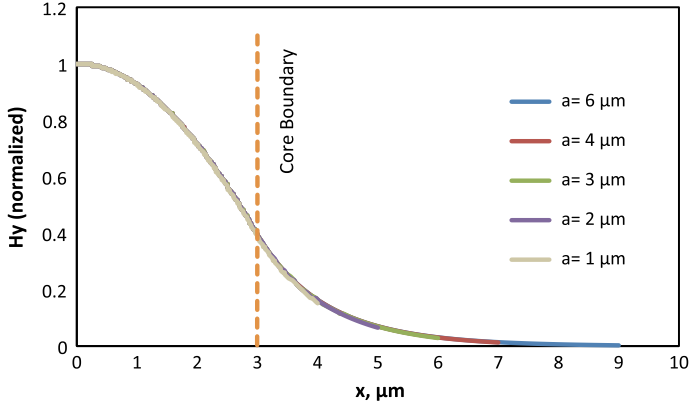
When infinite element is not used, the boundary wall acts like an electric wall. Hence, the effect of this wall is like a mirror, where  $H_y$  field is normal to this wall and forces a non-zero value which mimics an even coupled array. As even supermode of a coupled structure has higher effective index than that of isolated modes, so here the effective index becomes higher than that of the actual mode. Similarly, if the boundary on the top and bottom come closer to waveguide, this will force  $H_y = 0$  at the boundary and the effective index of  $H_{11}^y$  mode will reduce, as it mimics odd supermode. For the quasi-TM ( $H_{mn}^x$ ) mode, the boundary condition of the  $H_x$  field would be Neumann for the upper and lower boundary and Dirichlet for vertical side walls.

Figure 4 shows the  $H_y$  field variation along the x-direction with infinite element. Once again, the  $H_y$  field variation obtained so far for the different waveguide width is converted to a normalised form to make them comparable with each other. The plot shows that the field, even when the side wall was closer to waveguide, still follows the actual modal field with highest waveguide width. That means, even when the computational domain is reduced, it does not force the natural boundary condition at that position. Resulting fields correctly represent the field decay ( $e^{-\alpha x}$ ) in the cladding region.

Next, a high-index contrast waveguide structure is considered, where a Silicon core, with refractive index of 3.47638, is placed on a  $\text{SiO}_2$  buffer having refractive index of 1.44427 and covered with Air cladding. The core width is taken as 800 nm and its height as 200 nm. Using the **H**-field FEM (Rahman and Davies 1984), the numerical analysis has been carried out to find the effective index of the fundamental and second modes, for different mesh divisions. The structure considered here has an onefold symmetry. So, only



**Fig. 3** Effective index versus 'a' for Ge doped Silica

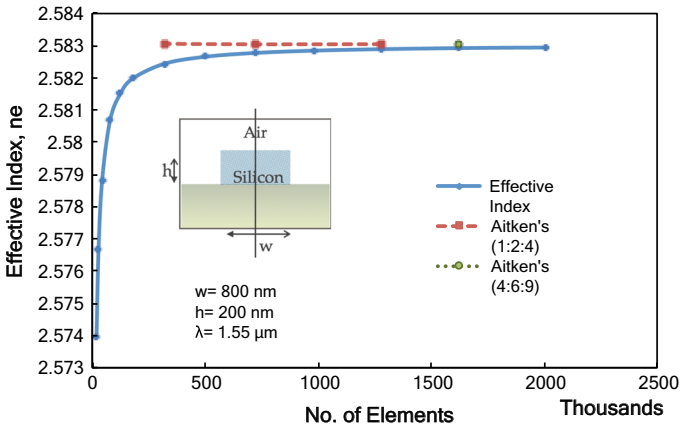


**Fig. 4**  $H_y$  field variation along width of waveguide for Ge doped Silica with infinite wall

half symmetry waveguide is represented and the simulation is conducted at 1550 nm operating wavelength.

The simulation result reveals that effective index increases with the finer mesh to a certain level and then settles asymptotically to 2.5829386, as shown by a blue line in Fig. 5. Similar as for low-index contrast, to test the convergence, Aitken extrapolation technique has been used, considering a fixed geometric ratio of 1:2:4 and 1:1.5:2.25 (integer 4:6:9). Thus, the new effective index is calculated using Eq. (9), shown by red dotted line for 1:2:4 and green dot for 4:6:9. Based on the results, it is observed that the convergence value is more accurate with the geometric ratio of 4:6:9 for a given final mesh division used. Although using a geometric ratio 4:6:9 is more restrictive in distributing mesh over the whole problem domain, but we have observed that this approach yields better convergence than using a simpler 1:2:4 geometric mesh refinement.

Table 2 shows the effective index of the fundamental quasi-TE  $H_{11}^y$  mode for a Silicon Nanowire. For comparison purpose, the commercially available packages, COMSOL, FIMMPROP and FDTD based Lumerical were used to find the effective index of the same



**Fig. 5** Effective index versus number of elements for silicon nanowire

**Table 2** Effective index of Si nanowire using various numerical methods

FEM (half structure)	Aitken's	COMSOL	FIMMPROP	Lumerical
2.5829386	2.5830376	2.58293381	2.58307	2.581929

waveguide. The results shows effective indices are nearly close to each other with FIMMPROP result slightly higher and that of Lumerical considerably lower, which are also tabulated in Table 2. It can be noted that for this case, when the index contrast is higher, variations of effective index with the mesh or between the approaches are higher than that of a low index contrast Silica guide, as shown in Table 1.

However, as the  $\Delta n$  between the core and cladding index was higher for a Silicon nanowire, it may be useful to compare the range of normalized propagation constant, **b**. The normalised propagation constant, 'b' variation with mesh size for Silicon Nanowire is also found to be significantly high with 0.0045. This suggests, extra care must be taken to find accurate solutions of high index contrast waveguides.

### 3.2 Directional coupler

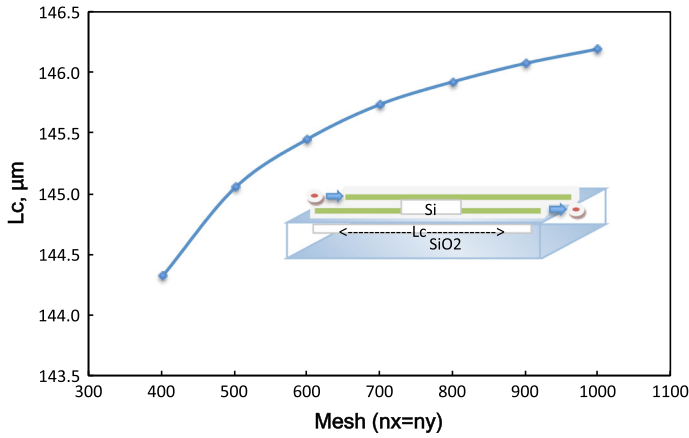
The directional coupler is a key optical devices where modal solution can also be applied, and yet z-dependence optical parameter extracted from different numerical approaches can be tested. It works on the principle that the two guides, having separation,  $s$ , and with light input into Guide 1 is completely coupled into Guide 2 when the length of the device,  $L_c$ , is,

$$L_c = \frac{\pi}{\beta_e - \beta_o} \quad (14)$$

where,  $\beta_e$  and  $\beta_o$  are the propagation constants of the even and odd supermodes.

Using **H**-field FEM, the simulation of Silicon waveguide directional coupler was performed considering different width, height and separation ranging between 100 and 1000 nm. For these dimensions, simulation has been conducted for  $400 \times 400$  to  $1000 \times 1000$  mesh divisions in order to analyse its trend. Typical result for the coupling length of the Silicon waveguide directional coupler for 800 nm core width, 200 nm core height and 100 nm separation for 1550 nm operating wavelength, with mesh division of  $1000 \times 1000$  is found to be 146.19  $\mu\text{m}$  that seems to converge with very little increment if the simulation is carried out beyond  $1000 \times 1000$  mesh. For the **H**-formulation used here, the individual propagation constants,  $\beta_e$  and  $\beta_o$  of the two supermodes increase slightly with the mesh, but their difference is more stable with the mesh division. The coupling length variation against mesh division is shown in Fig. 6, which shows its convergence as the mesh number is increased. However, their differences reduces slightly, so with the increasing mesh,  $L_c$  increases upto the saturation level.

Using commercial package, COMSOL, FIMMPROP and FDTD based Lumerical, the simulations were carried out for the Directional Coupler with the same physical dimension as above. Results show that the coupling length obtained from COMSOL and FIMMPROP were very close to the FEM results but slightly lower in case of Lumerical. The results are tabulated in Table 3.



**Fig. 6** Variation of the coupling length with the mesh division

**Table 3** Coupling length,  $L_c$ , using various numerical methods

FEM	COMSOL	FIMMPROP	Lumerical
146.19	146.1	144.8	142.85

## 4 Conclusion

The Finite Element Method and Finite-Difference Time-Domain Method are very popular today in terms of solving electromagnetic problems. However, new codes, packages or methods are constantly published with novel analytical methods to solve electromagnetic problems. For any photonics modelling, it is useful to have some confidence in the results obtained.

In this paper, we aimed to provide a detailed and comprehensive analysis defining the modal characteristics of both low-index contrast Silica and high-index contrast Silicon nanowire waveguides and these are presented here. We have also shown the effects of the mesh size used and the infinite element. The results obtained so far are again compared with various other numerical methods in order to benchmark them. It can be noted that, as the solution accuracy of high index guide is slightly poorer, convergence of the modal solutions should be checked and if possible they should be benchmarked against another alternative approaches.

It is also shown the advantage of exploiting the structural symmetry, if available, to obtain better accuracy and also to avoid mode degeneration. Similarly, more accurate solution can be obtained by using Aitken's extrapolation. Most of the commercial package do not clearly state the boundary conditions at the orthodox boundary, and it is shown here that if the boundary is taken close to waveguide core it can introduce error in both the effective index value and for the field profile. If possible, infinite element can be introduced to avoid artificial effect of the computational window. It should be noted that PML boundary (Berenger 1994) should only be used when leakage or bending loss needs to be

calculated, as introduction of PML makes the eigenvalue equation complex, and needs more computational resources.

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