Nanophotonic Computational Design

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Abstract: In contrast to designing nanophotonic devices by tuning a handful of device parameters, we have developed a computational method which utilizes the full parameter space to design linear nanophotonic devices. We show that our method may indeed be capable of designing any linear nanophotonic device by demonstrating designed structures which are fully three-dimensional and multi-modal, exhibit novel functionality, have very compact footprints, exhibit high efficiency, and are manufacturable. In addition, we also demonstrate the ability to produce structures which are strongly robust to wavelength and temperature shift, as well as fabrication error. Critically, we show that our method does not require the user to be a nanophotonic expert or to perform any manual tuning. Instead, we are able to design devices solely based on the user's desired performance specification for the device.

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References and links

1. Introduction

Currently, almost all nanophotonic components are designed by hand-tuning a small number of parameters (e.g. waveguide widths and gaps, hole and ring sizes). However, the realization of increasingly complex, dense, and robust on-chip optical networks will require utilizing increasing numbers of parameters when designing nanophotonic components.

Opening the design space to include many more parameters allows for smaller footprint, higher performance devices by definition; since original designs are still included in this parameter space. Unfortunately, the lack of intuition for what such designs might look like and the inability to manually search such a large parameter space have greatly hindered the ability to employ anything even close to the available parameter space for designing nanophotonic components.

For this reason, we have developed and implemented a computational method which is able to use the full parameter space to design linear nanophotonic components in three dimensions. Critically, our method requires no user intervention or manual tuning. Instead, a *design-by-specification* scheme is used to produce designs based solely on a user's performance specification.

We show that our method can indeed produce designs which are extremely compact, and, at the same time, highly efficient. Furthermore, we demonstrate that devices with novel functionality are easily designed. We also show that our method can be used to produce designs with extreme robustness to wavelength and temperature shift, as well as fabrication error. Lastly, since all our results are produced by simply specifying the functionality and performance of the desired device, our results suggest that our method may indeed by able to design *all* linear nanophotonic devices.

2. Problem formulation

In order to produce designs which utilize the full parameter space, and are based solely on the user's performance specification, we formulate the design problem in the following way:

minimize
$$\sum_{i}^{M} ||A_i(z)x_i - b_i||^2$$
 (1a)

subject to
$$\alpha_{ij} \le |c_{ij}^{\dagger} x_i| \le \beta_{ij}$$
, for $i = 1, ..., M$ and $j = 1, ..., N_i$ (1b)

$$z_{\min} \le z \le z_{\max}$$
 (1c)

The explanation for the various terms in eq. (1) follows:

1. $A_i(z)x_i - b_i$ is the *physics residual* for the *i*th mode. That is to say, $A_i(z)x_i - b_i$ represents the underlying physics of the problem; namely, the electromagnetic wave equation $(\nabla \times \mu_0^{-1} \nabla \times -\omega_i^2 \varepsilon) E_i + i\omega_i J_i$.

The specific substitutions used in order to transform

$$(\nabla \times \mu_0^{-1} \nabla \times -\omega_i^2 \varepsilon) E_i + i \omega_i J_i \longrightarrow A_i(z) x_i - b_i$$

are

- $E_i \rightarrow x_i$,
- $\varepsilon \rightarrow z$,
- $\nabla \times \mu_0^{-1} \nabla \times -\omega_i^2 \varepsilon \to A_i(z)$, and
- $-i\omega_i J_i \rightarrow b_i$.

In contrast to typical schemes for optimizing physical structures, our formulation actually allows for non-zero physics residuals; which can be deduced since $A_i(z)x_i - b_i = 0$ is not a hard constraint. Instead, this formulation is what we call an *objective-first* formulation in that the *design objective* (explained below) is prioritized above satisfying physics.

2. The (field) design objective consist of the constraint $\alpha_{ij} \leq |c_{ij}^{\dagger}x_i| \leq \beta_{ij}$. Physically, this constraint describes the performance specification of the device via a series of field overlap integrals at various output ports of the device. Specifically, the $c_{ij}^{\dagger}x_i$ terms represents an overlap integral between the E-field of the *i*th mode (x_i) with an E-field of the user's choice (c_{ij}) , where the additional subscript j allows the user to include multiple such fields. The amplitude of the overlap integral is then forced to reside between α_{ij} and β_{ij} .

This mechanism allows the user to express the desired performance of the device as a combination of field amplitudes in various output field patterns. These outputs would be in response to a predefined input excitation, which is determined by the current excitation b_i ($-i\omega_i J_i$) in the physics residual of each mode.

As an example of a design objective for some mode 1 a user might choose to have the majority of the output power reside in some output pattern 1, while ensuring that only a small amount of power be transferred to some output pattern 2. In this case the user

would use $0.9 \le |c_{11}^\dagger x_1| \le 1.0$ for the former, and then $0.0 \le |c_{12}^\dagger x_1| \le 0.01$ for the latter; where c_{11} and c_{12} are representative of output patterns 1 and 2 respectively.

Finally, we note again that the design objective in our formulation is actually a hard constraint. This means that it is *always satisfied*, even to the extent of allowing for an unphysical field (since the physics residual will not be exactly 0). It is for this reason that we call such a formulation "objective-first".

3. The final term in eq. (1), $z_{\min} \le z \le z_{\max}$, is the structure design objective. It is used as a relaxation of the binary constraint, $z \in z_{\min}, z_{\max}$, which would ensure that the final design be composed of two discrete materials.

3. Method of solution

We employed the alternating directions method of multipliers (ADMM) algorithm in order to solve eq. (1). The ADMM algorithm solves eq. (1) by iteratively solving for x_i , z, and a dual variable u_i .

Since we are working in three dimensions, solving eq. (1) for x_i is non-trivial in that it involves millions of variables and requires solving for the ill-conditioned $A_i(z)$ matrix. For this reason, we use a home-built finite-difference frequency-domain (FDFD) solver which implements a hardware-accelerated iterative solver on Amazon's Elastic Compute Cloud. Critically, our cloud-based solver allows us to scale to solve problems with arbitrarily-large number of modes, with no significant penalty in runtime.

In contrast to solving for x_i , solving for z is much simpler since we only consider planar structures; thereby limiting z to have only thousands of variables.

Lastly, in order to arrive at fully discrete, manufacturable structures, we convert z to a boundary parameterization and tune our structure using a steepest-descent method.

4. Results

5. Conclusion