

Proof-of-concept implementation of the massively parallel algorithm for simulation of dispersion-managed WDM optical fiber systems

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We perform a proof-of-concept implementation of the massively parallel algorithm [P. M. Lushnikov, Opt. Lett. 27, 939 (2002)] for simulation of dispersion-managed wavelength-division-multiplexed optical fiber systems. Linear scalability of the algorithm with the number of computer cores is demonstrated. Exact result on the accuracy of the implemented algorithm is found analytically and confirmed numerically as well as it is compared with the accuracy of the standard split-step algorithm. © 2011 Optical Society of America

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A wavelength-division-multiplexed (WDM) dispersion-managed (DM) optical fiber system is the basis of current high-bit-rate optical communications. The next generation of these systems will use both the amplitude and phase of the optical signal as a carrier of information (see, e.g., [1,2]) to achieve higher system performance. WDM systems are weakly nonlinear ones with a linear dispersion length typically in the range of tens of kilometers while a nonlinear length is at several hundreds of kilometers [3–6]. Nonlinearity is a major factor limiting performance of such systems, while linear effects can be significantly compensated by coherent detection.

WDM requires propagation of a wide range of frequencies through optical fiber coupled by the nonlinearity. Path-averaged group-velocity dispersion GVD effects cause optical pulses in distinct WDM channels to move with different group velocities. Consequently, modeling of WDM systems requires simulating a long time interval, which determines needed high resolution in the frequencies. All that makes accurate numerical simulations enormously challenging with very large number of Fourier modes N needing to be resolved. The standard algorithm for such simulation is an operator splitting, or split-step algorithm (SS). It involves several Fourier transforms for every spatial step along an optical fiber. The fast Fourier transform (FFT) algorithm computes each such transformation in $O(N \log N)$ operations of multiplication. The efficiency of using supercomputers for SS is limited because parallel algorithms for one-dimensional FFT (contrary to multidimensional FFT) provide only very moderate speed up. For example, one of the leading implementations [7] shows at best four times acceleration on 16 processor cores on the system with shared memory. Further increase of the CPU number appears to be inefficient. Further increase of the number of processors on conventional systems requires use of a distributed memory approach (cluster), which has higher latency of nodes interconnection media. The experimental data shows [7] that at a number of harmonics up to 2^{21} , a shared memory approach is more efficient. At a higher number of harmonics, moderate

acceleration can be achieved on a cluster, although scaling would still be far from linear.

Here we demonstrate the proof-of-concept realization of the massively parallel algorithm (MPA) for simulation of WDM systems that is free from all these limitations. MPA was proposed by one of the authors of this Letter in [8] and exploits weak nonlinearity of the WDM system. We demonstrate the linear scalability of performance with number of computer cores. We also obtain the exact result on the accuracy of the algorithm in comparison with SS. The results are in full agreement with numerics.

We neglect polarization effects, stimulated Raman scattering, and Brillouin scattering. Then the propagation of WDM optical pulses in DM fiber is described by a scalar nonlinear Schrödinger equation

$$iA_z - \frac{1}{2}\beta_2(z)A_{tt} - \frac{i}{6}\beta_3(z)A_{ttt} + \sigma(z)|A|^2A = iG(z)A, \quad (1)$$

where $G(z) \equiv \{-\gamma + [\exp(z_a\gamma) - 1]\sum_{k=1}^N \delta(z - z_k)\}$; z is the propagation distance along an optical fiber; $A(t, z)$ is the slow amplitude of light; β_2 and β_3 are the first- and second-order GVD, respectively, which are periodic functions of z ; $\sigma = (2\pi n_2)/(\lambda_0 A_{\text{eff}})$ is the nonlinear coefficient; n_2 is the nonlinear refractive index; λ_0 is the carrier wavelength; A_{eff} is the effective fiber area; $z_k = kz_a$ ($k = 1, 2, \dots, N$) are amplifier locations; and γ is the loss coefficient. Distributed amplification can be also included in $G(z)$.

Applying Fourier transform $\hat{A}(\omega, z) = \hat{F}[A(t, z)] = \int_{-\infty}^{\infty} A(t, z) \exp[i\omega t] dt$ to Eq. (1), using the change of variables $\hat{A}(\omega, z) \equiv \hat{\psi}(\omega, z) \exp[i\beta(\omega, z) + \int_{z_0}^z G(z') dz']$ and integration over z results in the integral equation

$$\hat{\psi}(\omega, z) = \hat{\psi}(\omega, z_0) + i \int_{z_0}^z \sigma(z') \hat{F}[|A(t, z')|^2 A(t, z')] \times e^{-i\beta(\omega, z') - \int_{z_0}^{z'} G(z'') dz''} dz', \quad (2)$$

where $\beta(\omega, z) = \int_{z_0}^z [\frac{\omega^2}{2}\beta_2(z') + \frac{\omega^3}{6}\beta_3(z')] dz'$.

Case $\hat{\psi}(\omega, z) = \text{const}$ corresponds to the exact solution of the linear part of Eq. (1) [or, equivalently, setting $\hat{F}[\cdot] \equiv 0$ in (2)]. Assume that the nonlinearity is weak, $z_{\text{nl}} \gg z_{\text{disp}}$, where $z_{\text{nl}} \equiv 1/|p|^2$ is a characteristic nonlinear length, $z_{\text{disp}} \equiv \tau^2/|\beta_2|$ is the dispersion length, and p and τ are typical pulse amplitude and width, respectively. Then $\hat{\psi}(\omega, z)$ is a slow function of z on any scale $L \ll z_{\text{nl}}$ (see [4,9,10]). We solve Eq. (1) by iterations for $0 \leq z - z_0 \leq L$, where we have a freedom of choice of L with the only condition that $L \lesssim z_{\text{disp}}$. For the first iteration we set $\hat{\psi}^{(0)}(\omega, z) = \hat{\psi}(\omega, z_0) = \hat{A}(\omega, z_0)$ and, respectively, $A(z, \omega) = \hat{\psi}^{(0)}(\omega, z) e^{i\beta(\omega, z) + \int_{z_0}^z G(z') dz'}$ on the right-hand side (rhs) of Eq. (1), which gives the first iteration $\hat{\psi}^{(1)}(\omega, z)$ for the left-hand side (lhs) of Eq. (1). Similarly, substitution of $\hat{\psi}^{(n-1)}(\omega, z)$ in the rhs of Eq. (1) gives $\hat{\psi}^{(n)}(\omega, z)$ in the lhs of Eq. (1) for $n = 1, 2, \dots$.

In simulations we use $\hat{\psi}(\omega, z_0)$ with $z_0 = mL$ for a given $m = 0, 1, \dots$ to perform a total number of iterations n_{tot} to approximate $\hat{\psi}(\omega, z_0 + L)$ as $\hat{\psi}^{(n_{\text{tot}})}(\omega, z_0 + L)$. Then we use that approximate value as starter for the next spatial interval by setting $z_0 = (m + 1)L$ and proceeding in a similar way.

Assume that the interval $z_0 \leq z \leq z_0 + L$ includes $M + 1$ equally spaced points $z_0, z_1, \dots, z_M = z_0 + L$. The MPA is based on these iterations as follows:

1. $\hat{A}(\omega, z_0) = \hat{\psi}(\omega, z_0)$, copy $\hat{\psi}(\omega, z_0)$ in $\hat{\psi}(\omega, z)$ at all z .
2. Find $\hat{A}(\omega, z) = \hat{\psi}(\omega, z) e^{i\beta(\omega, z) + \int_{z_0}^z G(z') dz'}$ at all z .
3. In order to return to t -domain, calculate independent Fourier transforms $A(t, z) = \hat{F}^{-1}[\hat{A}(\omega, z)]$ at all z .
4. Calculate independent Fourier transforms $\hat{V}(\omega, z) = \hat{F}[|A(t, z)|^2 A(t, z)]$ at all z .
5. Numerical integration (summation) by trapezoidal rule of the integral in Eq. (2) using $\hat{V}(\omega, z)$ from step 4. Save intermediate results of integration at every z .
6. For the second, third, etc., iterations go to step 2.
7. Reconstruct $\hat{A}(\omega, z_M)$ on the far edge of interval.

The MPA is schematically shown in Fig. 1. All steps of the MPA are computed parallelly. The most time-consuming steps are 3 and 4. All FFTs at each z are independently performed in CPU cores (vertical bars in Fig. 1). Calculations in steps 2 and 5 are done for every harmonic independently (dashed horizontal arrow in Fig. 1).

We implement the MPA for shared memory symmetric multiprocessor (SMP) architecture. The only powerful SMP computer in exclusive use was the HP SuperDome

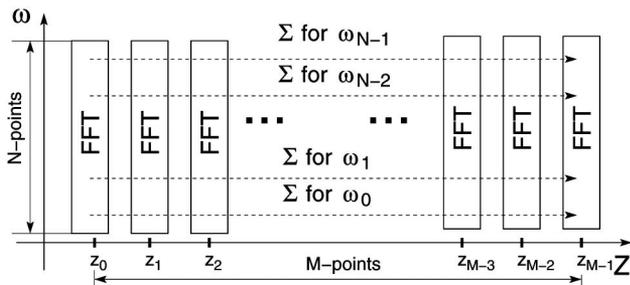


Fig. 1. Schematic representation of MPA.

64000 supercomputer, equipped with 64 HP PA-RISC processors (<http://jscc.ru>). Both processors and memory bandwidth are outdated and relatively slow. However, for the proof-of-concept simulation the main criteria is the number of processors in the system.

Simulations were performed in a setup identical to the one used in [8] with pseudorandom sequences of optical pulses in five channels of 20 periods of a WDM DM system. The main difference between the current MPA implementation and the original algorithm [8] is in a more efficient way to handle summation in step 5, optimizing CPUs cache use. We achieve ~ 30 times speed up with respect to a single processor version of the code. Figure 2 shows the scalability of performance. Scaling is close to linear up to 32 processors. Then memory bandwidth limitation of the available SMP computer makes further parallelization less efficient. Another reason for that was restriction on the memory usage, which limited the number of Fourier harmonics. However, we see a clear tendency of the scalability improvement with an increase of the number of harmonics because the longer time the computer spends in FFTs the less important communications are.

To find the accuracy of the MPA we put the exact solution of Eq. (1) in the operator form as

$$A_{\text{exact}}(z) = \exp[i(\hat{L} + \hat{N})z]A(0), \quad (3)$$

where \hat{L} represents all linear terms and \hat{N} represents the nonlinear term in Eq. (1), and we set $z_0 = 0$. Here and below for brevity we omit the argument t of function $A(z, t)$. We assume that $\beta_2(z)$, $\beta_3(z)$, $\sigma(z)$, $G(z)$ are constant functions of z at each interval of length L , although a generalization to more general dependence on z is straightforward. SS uses the efficiency and high precision of the simulations for $\exp[i\hat{L}z]A(0)$ and $\exp[i\hat{N}z]A(0)$. But \hat{L} and \hat{N} do not commute and we need to approximate Eq. (3) for $z = L$ by the composite M steps of SS as

$$A_{\text{SS}}(L) \equiv \exp[i\hat{L}\Delta z/2]\hat{Q}^M \exp[-i\hat{L}\Delta z/2]A(0), \quad (4)$$

where $\Delta z \equiv L/M$, $\hat{Q} \equiv \exp[i\hat{N}\Delta z]\exp[i\hat{L}\Delta z]$. Taylor series expansion of operators in Eqs. (4) and (3) gives the following error $r_1 \equiv A_{\text{exact}}(L) - A_{\text{SS}}(L)$ of composite SS for arbitrary M assuming $L \lesssim z_{\text{disp}} \ll z_{\text{nl}}$:

$$r_1 = i[L^3/(2M^2)]P_L + [L^3/M^2]O(\hat{L}\hat{N}^2A_0), \quad (5)$$

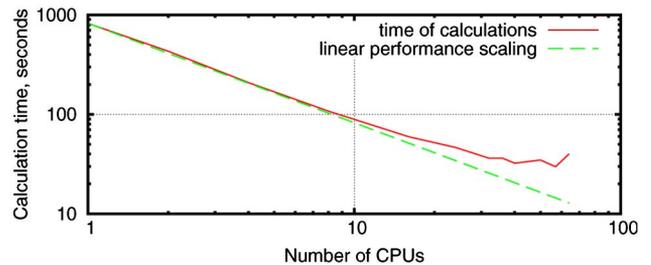


Fig. 2. (Color online) Scalability of MPA on HP SuperDome 64000.

where $A_0 \equiv A(0)$, $O(\hat{\mathcal{L}}^k \hat{\mathcal{N}}^l A_0)$ means different combinations of terms with k th power of $\hat{\mathcal{L}}$ and l th power of $\hat{\mathcal{N}}$. Also, $P_L = \frac{1}{6} \bar{A}_0 (\hat{\mathcal{L}} A_0)^2 - \frac{1}{3} A_0 |\hat{\mathcal{L}} A_0|^2 + \frac{1}{6} |A_0|^2 \hat{\mathcal{L}}^2 A_0 - \frac{1}{3} \hat{\mathcal{L}} (|A_0|^2 \hat{\mathcal{L}} A_0) + \frac{1}{12} A_0^2 \hat{\mathcal{L}}^2 \bar{A}_0 + \frac{1}{6} \hat{\mathcal{L}} (A_0^2 \hat{\mathcal{L}} \bar{A}_0) + \frac{1}{12} \hat{\mathcal{L}}^2 (|A_0|^2 A_0)$ represents all terms with the second power in $\hat{\mathcal{L}}$. Note that the operator expansions for arbitrary M are not trivial and require the extensive use of the symbolic computations.

Discretization of iterations over z in Eq. (2) with $\hat{\psi}_j(\omega) \equiv \hat{\psi}(z_j, \omega)$, $z_j = j\Delta z$, $j = 0, 1, \dots, M$ at each N discrete values of ω is given by

$$\hat{\psi}_l^{(n+1)}(\omega) = \hat{\psi}_0^{(n)}(\omega) + i \frac{V_0^{(n)}}{2} + i \sum_{j=1}^{l-1} V_j^{(n)} + i \frac{V_l^{(n)}}{2}, \quad (6)$$

where $l = 1, 2, \dots, M$, $V_j^{(n)} \equiv \sigma(z_j) \hat{F}[|A_j^{(n)}|^2 A_j^{(n)}] \Delta z \exp[-i\beta(\omega, z_j) - \int_0^{z_j} G(z) dz]$, $\hat{A}_j^{(n)} = \hat{\psi}_j^{(n)} \exp[i\beta(\omega, z_j) + \int_0^{z_j} G(z) dz]$, and $\hat{\psi}_l^{(n)}(\omega)$ is the n th iteration of ψ while for zero iteration $\hat{\psi}_j^{(0)} = A_0$, $j = 0, 1, 2, \dots, M$. From comparison of the operator expansion for n th iteration with the operator expansion of the exact solution in Eq. (3) we obtain (again assuming that $\beta_2(z)$, $\beta_3(z)$, $\sigma(z)$, $G(z)$ are constant functions of z) that the error $r_2 \equiv A_{\text{exact}}(L) - A^{(n)}(L)$ of composite SS for arbitrary M :

$$r_2 = i \frac{L^3}{M^2} P_L + \frac{L^3}{M^2} O(\hat{\mathcal{L}} \hat{\mathcal{N}}^2 A_0) + O(\hat{\mathcal{N}}^{n+1} A_0), \quad (7)$$

where P_L is the same as in Eq. (5) and we assume that $n \geq 3$, which ensures that r_2 at leading order $O(L^3)$ does not depend on n . For $n = 2$ the additional error term is $O([L^4/M^2] \hat{\mathcal{L}}^3 \hat{\mathcal{N}} A_0)$, which can be of the same order as $L^3 P_L / M^2$ provided $L \sim z_{\text{disp}}$. But for practical realization of the MPA we expect that $L \ll z_{\text{disp}}$, and then $n_{\text{tot}} = 2$ can be also an optimal choice.

Error term αP_L dominates in Eq. (5), while in Eq. (7) it competes with the last term on the rhs, which has the order $O(\hat{\mathcal{N}}^{n+1} A_0) \sim (L/z_{\text{nl}})^{n+1} A_0$ and is independent of M because it results from the iterations of Eq. (2) in the continuous limit $M \rightarrow \infty$. An increase of n ensures dominance of αP_L in Eq. (7) because $L \lesssim z_{\text{disp}} \ll z_{\text{nl}}$. Then we conclude from comparison of Eqs. (7) and (5) that SS error is twice smaller than the MPA. So to match the accuracy of SS it is enough for the MPA to take M by a factor $2^{1/2}$ larger. Respectively, the MPA requires a minimum $2^{1/2}n + 1$ CPU cores to outperform SS ($2^{1/2}n$ would be the exact match of performance). For example, in simulations with the parameters of Fig. 2, 250 DM periods and $n = 2$, we obtained the ratio 9.2 of SS and MPA computation times at equal accuracy and 32 cores, which is close to the theoretical prediction $2^{-1/2} 32 n^{-1}$.

To check these analytical predictions we simulated a three-channel WDM system over one period of a DM fiber system with 2^{11} frequency harmonics, $L = 20$ km for the standard fiber, and other parameters as in [8]. As a

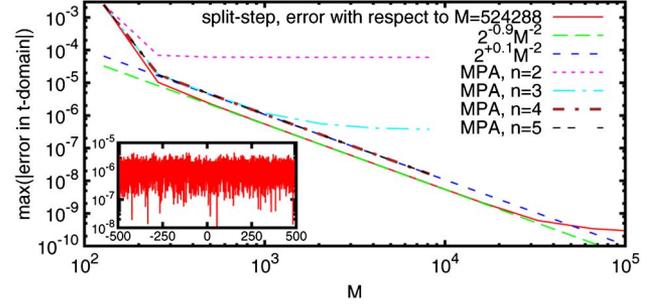


Fig. 3. (Color online) Errors in L_∞ norm (maximum over t) of MPA and SS versus M . Results for fourth and fifth iterations are visually indistinguishable with the theoretical scaling line. Inset shows the error (normalized to $\max|\psi(t)|$) with t for 20-channel WDM system after 10^4 km.

“numerically exact” we used SS with 2^{21} grid steps over one DM system period. Figure 3 shows that the error of the MPA with $n = 3$ scales as one for SS. That is, $n = 3$ is enough to neglect $O(\hat{\mathcal{N}}^{n+1} A_0)$ term in Eq. (7). Errors for MPA and SS are different by a factor 2 in full agreement with Eqs. (5) and (7). We also compared MPA and SS for simulation of the transoceanic distance 10^4 km (250 DM periods) of the realistic WDM system with 20 channels using $N = 2^{13}$, $M = 2^{14}$, $L = 1.25$ km. That system has $(20/3)^2$ higher nonlinearity than above so we decreased L . The inset of Fig. 3 shows the error of the MPA with $n = 3$ in that case. A ratio of MPA and SS L_∞ errors (i.e., max over t in that inset and similar for SS) is $\simeq 2.1$, again close to 2.

In conclusion, we have demonstrated the feasibility of the MPA. Scaling of the parallel version on the available SMP machine was close to linear up to 32 processing threads even on outdated architecture and with a very restricted size of FFT arrays. We propose using a shared memory model for parallel computation, which has lower penalties due to interprocess communications, and exploiting the power of modern graphics processing units (GPUs). Nvidia Tesla C2070 GPU has 448 cores and 6 GB memory, which appears quite suitable for the MPA.

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