

# Mixed order discretization based two-level Schwarz preconditioners for a tracer transport problem on the cubed-sphere



Haijian Yang<sup>a</sup>, Chao Yang<sup>b,c,\*</sup>, Xiao-Chuan Cai<sup>d</sup>

<sup>a</sup> College of Mathematics and Econometrics, Hunan University, Changsha, Hunan 410082, China

<sup>b</sup> Institute of Software, Chinese Academy of Sciences, Beijing 100190, China

<sup>c</sup> State Key Laboratory of Computer Science, Chinese Academy of Sciences, Beijing 100190, China

<sup>d</sup> Department of Computer Science, University of Colorado Boulder, Boulder, CO 80309, USA

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## ABSTRACT

Solving tracer transport problems on the sphere at high resolution is important for atmospheric science. In this paper, we study a fully implicit method that allows larger time step size than explicit methods. The key to the success of the implicit approach is how to solve the discretized linear system at each time step. We investigate a parallel Krylov–Schwarz approach based on two discretizations, and test the proposed algorithm with two benchmark test cases. We show numerically that the solver scales well in terms of both strong and weak scalabilities on a supercomputer with thousands of processors.

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## 1. Introduction

The tracer transport problem plays an important role in global atmospheric modeling [5]. Its numerical simulation at high resolution is very demanding in terms of computational resources. In recent years, there have been numerous efforts to develop advanced numerical methods for solving global atmospheric problems [7,9–11,14,20,21]. Among these, the fully implicit method has become an interesting research focus because the time step size is no longer constrained by the stability condition. The main challenge to implicit methods is how to solve the large system of equations.

In this paper, we study a Krylov–Schwarz method [2,12] to efficiently solve the linear system arising from the implicit finite volume discretization of a tracer transport problem. We use a class of two-level restricted Schwarz methods based on mixed order discretizations, in which the coarse preconditioner is based on a first-order spatial discretization while a second-order scheme is used to construct the preconditioner on the fine mesh. Our experiments show that this strategy leads to better parallel performance than using the second-order discretization on both levels.

The rest of the paper is organized as follows. In Section 2, we present a fully implicit discretization scheme for the tracer transport problem on the cubed-sphere. Section 3 focuses on the details of the two-level Schwarz preconditioning method. Numerical results for two benchmark test cases are provided in Section 4 to illustrate the accuracy and the parallel performance of the proposed methods. We end the paper with some concluding remarks in Section 5.

## 2. Finite volume discretization of the tracer transport problem

Let  $\mathbb{S}$  be the surface of the sphere that is a closed domain without boundary and  $T$  denote the final time of the simulation. In the absence of source or sink, we write the tracer transport equation [5] defined on  $\mathbb{S} \times (0, T]$  as

$$\frac{\partial \phi}{\partial t} + \nabla \cdot (\mathbf{V}\phi) = 0, \quad (1)$$

where  $\phi$  is the transported scalar,  $\mathbf{V}$  is a given two-dimensional velocity vector, and  $\nabla$  is the gradient operator. We discretize the surface of the sphere using the cubed-sphere mesh, which is mapped to the sphere from the six faces of an inscribed cube  $\Omega$  based on a gnomonic projection [9,10,14]. Let  $(\xi, \eta) \in [-\pi/4, \pi/4] \times [-\pi/4, \pi/4]$  be the curvilinear coordinates on the cubed-sphere. Then (1) defined on each of the six faces of  $\Omega$  can be written as

\* Corresponding author at: Institute of Software, Chinese Academy of Sciences, Beijing 100190, China. Tel./fax: +86 10 62661632.

E-mail addresses: [haijianyang@gmail.com](mailto:haijianyang@gmail.com) (H. Yang), [yangchao@iscas.ac.cn](mailto:yangchao@iscas.ac.cn) (C. Yang), [cai@cs.colorado.edu](mailto:cai@cs.colorado.edu) (X.-C. Cai).

$$\frac{\partial \phi}{\partial t} + \frac{1}{\Lambda} \left( \frac{\partial}{\partial \xi} (\Lambda v^1 \phi) \right) + \frac{1}{\Lambda} \left( \frac{\partial}{\partial \eta} (\Lambda v^2 \phi) \right) = 0, \tag{2}$$

where  $\Lambda = (\sec \xi \sec \eta)^2 / \sqrt{(1 + \tan^2 \xi + \tan^2 \eta)^3}$  and  $(v^1, v^2)$  is the contravariant coordinates of  $\mathbf{V}$ ; more details can be found in [10,13].

In the study, a cell-centered finite volume method is used to discretize (2) in the spatial domain  $\Omega$ , which consists of the six faces  $\Omega^k$  ( $k = 1, 2, \dots, 6$ ). Let each  $\Omega^k$  be covered with  $N \times N$  mesh cells. Then each cell  $\Omega_{ij}^k = [\xi_{i-1/2}, \xi_{i+1/2}] \times [\eta_{j-1/2}, \eta_{j+1/2}]$  is centered at the position  $\xi_i = -\pi/4 + ih$  and  $\eta_j = -\pi/4 + jh$  with  $i, j = 1, \dots, N$ , and  $h = \pi/(2N)$ . The approximate solution  $Q_{ij}^k$  in the cell  $\Omega_{ij}^k$  is defined as

$$Q_{ij}^k = \frac{1}{h^2 \Lambda_{ij}} \int_{\eta_{i-1/2}}^{\eta_{i+1/2}} \int_{\xi_{i-1/2}}^{\xi_{i+1/2}} \Lambda(\xi, \eta) \phi(\xi, \eta, t) d\Omega_{ij}^k,$$

where  $\Lambda_{ij}$  is computed at the cell center on each face. We integrate (2) over  $\Omega_{ij}^k$  and obtain the following semi-discrete system:

$$\frac{\partial Q_{ij}^k}{\partial t} + \frac{F_{i+1/2,j} - F_{i-1/2,j}}{\Lambda_{ij}h} + \frac{G_{ij+1/2} - G_{ij-1/2}}{\Lambda_{ij}h} = 0,$$

where the numerical fluxes  $F_{i\pm 1/2,j}$  and  $G_{ij\pm 1/2}$  defined on the four cell boundaries are approximated as

$$F_{i\pm 1/2,j} = \frac{\Lambda_{i\pm 1/2,j}}{h} \int_{\eta_{j+1/2}}^{\eta_{j-1/2}} v^1(\xi_{i\pm 1/2}, \eta) \phi(\xi_{i\pm 1/2}, \eta) d\eta,$$

$$G_{ij\pm 1/2} = \frac{\Lambda_{ij\pm 1/2}}{h} \int_{\xi_{i+1/2}}^{\xi_{i-1/2}} v^2(x, \eta_{j\pm 1/2}) \phi(x, \eta_{j\pm 1/2}) d\xi,$$

respectively. A local Lax–Friedrichs flux formula is employed for approximating  $F_{i\pm 1/2,j}$  and  $G_{ij\pm 1/2}$ :

$$\tilde{f}(Q^-, Q^+) = [(f(Q^-) + f(Q^+)) - \alpha(Q^+ - Q^-)]/2,$$

where  $\alpha$  is the maximum absolute value of the normal velocity along each cell boundary, and  $Q^-$  and  $Q^+$  are the reconstructed states of  $Q$  on the cell boundary.

Two approaches are implemented in our methods for the reconstruction of the left and right states. The first one uses a piecewise constant method by

$$\begin{cases} Q_{i-1/2,j}^- = Q_{i-1,j}, \\ Q_{i-1/2,j}^+ = Q_{i,j}, \\ Q_{i+1/2,j}^- = Q_{i,j}, \\ Q_{i+1/2,j}^+ = Q_{i+1,j}, \end{cases} \tag{3}$$

which leads to a first-order method [21]. To achieve a higher order of accuracy, we use a second-order reconstruction by

$$\begin{cases} Q_{i-1/2,j}^- = Q_{i-1,j} + (Q_{i,j} - Q_{i-2,j})/4, \\ Q_{i-1/2,j}^+ = Q_{i,j} - (Q_{i+1,j} - Q_{i-1,j})/4, \\ Q_{i+1/2,j}^- = Q_{i,j} + (Q_{i+1,j} - Q_{i-1,j})/4, \\ Q_{i+1/2,j}^+ = Q_{i+1,j} - (Q_{i+2,j} - Q_{i,j})/4. \end{cases} \tag{4}$$

We can similarly obtain  $Q^-$  and  $Q^+$  in the  $\eta$ -direction. On the cubed-sphere mesh, values near face boundaries should be correctly passed to couple the patches together. One layer of ghost cells is used and the numerical fluxes are calculated symmetrically across the interface [10,11,19,23].

### 3. Fully implicit discretization and domain decomposition preconditioning

A semi-discrete system is obtained after applying the spatial discretization to (2),

$$\frac{\partial Q_{ij}^k}{\partial t} = L(Q_{ij}^k). \tag{5}$$

To integrate (5) in time, we employ a second-order backward differentiation formula with a fixed time step size  $\Delta t$  and obtain

$$\frac{3Q_{ij}^{k,m} - 4Q_{ij}^{k,m-1} + Q_{ij}^{k,m-2}}{2\Delta t} = L(Q_{ij}^{k,m}), \tag{6}$$

where  $Q_{ij}^{k,m}$  is the numerical solution at the  $m^{\text{th}}$  time step. A first-order backward Euler method is applied at the first time step.

A major advantage of fully implicit schemes is that the time step size  $\Delta t$  is no longer constrained by the Courant–Friedrichs–Lewy (CFL) condition, which is often required by explicit or semi-implicit methods. But the price to pay in using implicit methods for the tracer transport problem is to solve a linear system (6) at each time step. In this study we use a Krylov–Schwarz method to fix this issue. More precisely speaking, we use a right-preconditioned Krylov subspace method (GMRES) [12] as the linear solver, and an overlapping one-level or two-level Schwarz method [2,4,18,22] as the preconditioner. The rest of the paper will focus on the use of one-level and two-level overlapping Schwarz preconditioners.

To define an overlapping restricted additive Schwarz preconditioner, we first partition the computational domain  $\Omega$  into  $N_p$  non-overlapping subdomains  $\Omega_l$  for  $l = 1, 2, \dots, N_p$ , such that

$$\overline{\Omega} = \bigcup_{l=1}^{N_p} \overline{\Omega}_l,$$

where  $N_p$  is the number of subdomains and also the number of processor cores. Then within  $\Omega$  we extend each subdomain  $\Omega_l$  with  $\delta$  layers of mesh cells to a larger subdomain  $\Omega_l^\delta$  that overlaps with its neighbors. In each overlapping subdomain, we define a local subdomain matrix  $A_l$  that is the restriction of the global matrix  $A$  to  $\Omega_l^\delta$  with the restriction operator  $R_l^\delta$ , i.e.,

$$A_l = R_l^\delta A (R_l^\delta)^T.$$

Here,  $R_l^\delta$  and  $(R_l^\delta)^T$  are the restriction and prolongation operators, respectively. Given a global vector defined on all mesh cells in  $\Omega$ ,  $R_l^\delta$  restricts the global vector to a local vector that is defined only on the mesh cells within the overlapping subdomain  $\Omega_l^\delta$ , while  $(R_l^\delta)^T$  prolongates the restricted vector back to a vector defined on the whole domain  $\Omega$  with zeros filled to the unknowns corresponding to mesh cells outside  $\Omega_l^\delta$ . The one-level restricted additive Schwarz (RAS) preconditioner is defined as

$$M_h^{-1} = \sum_{l=1}^{N_p} (R_l^\delta)^T (A_l)^{-1} R_l^\delta, \tag{7}$$

where the restriction operator  $R_l^\delta$  is defined as the restriction to the non-overlapping subdomain  $\Omega_l$ , and the matrix–vector multiplication with  $(A_l)^{-1}$  is obtained approximately by an incomplete LU (ILU) factorization.

Next we define a two-level Schwarz preconditioner by combining the one-level additive Schwarz preconditioner  $M_h^{-1}$  on the fine mesh with a coarse mesh preconditioner  $M_H^{-1}$  in a multiplicative manner. Let  $A_h$  ( $A_H$ ) be the global matrix on the fine (coarse) mesh and  $\mathcal{I}_H^h$  an interpolation operator from the coarse mesh to the fine mesh. Similarly,  $\mathcal{I}_h^H$  denotes a restriction operator from the fine mesh to the coarse mesh. Then a multiplicative type two-level Schwarz preconditioner [15,16,18] is defined as

$$M_{\text{two}}^{-1} = B_H^{-1} + M_h^{-1} - M_h^{-1} A_h B_H^{-1}, \tag{8}$$

where  $B_H^{-1} = \mathcal{I}_H^h A_H^{-1} \mathcal{I}_h^H$ . More precisely speaking, the application of the preconditioner  $y = M_{\text{two}}^{-1} r$  is obtained in the following steps:

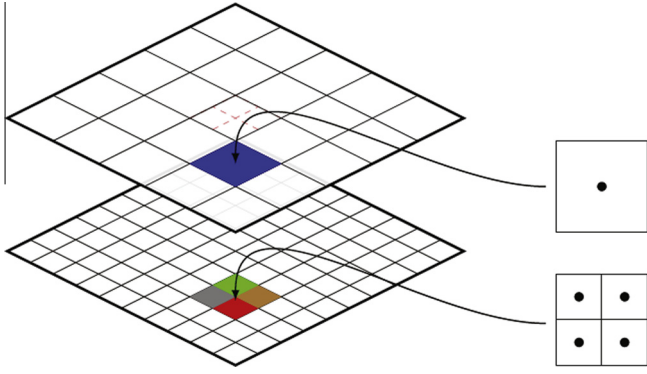


Fig. 1. Partitions for one face of the cubed-sphere with the coarse-to-fine mesh ratio 1:2. The bottom (upper) panel is the partition of the fine (coarse) mesh.

$$y = u_h + M_h^{-1}(r - A_h u_h). \tag{12}$$

In (10), a smaller linear system associated with  $A_H$  is then solved. Here, we solve the coarse problem iteratively by using GMRES with the one-level RAS preconditioner (7), i.e.,  $u_H = A_H^{-1}r_H$  is computed by approximately solving the following problem

$$\|r_H - A_H M_H^{-1} M_H u_H\| \leq \eta_r^c \|r_H\|,$$

where  $\eta_r^c$  is the solver tolerance for the iteration on the coarse mesh and  $M_H^{-1}$  is the one-level RAS preconditioner (7) on the coarse mesh.

On the coarse level, (10) is expensive to solve when the mesh size is fine. The situation becomes worse when a high order discretization is used, because the resulting matrix has considerably more non-zero elements. On the other hand, the coarse solve is used to improve the convergence of the iterative solver, it may not need to be solved as accurately as on the fine mesh. Based on these considerations, we propose a mixed order discretization based two-level Schwarz preconditioner. The approach consists of two components: we solve the coarse problem based on the sub-domain matrix with the first-order spatial discretization (3) for the construction of the preconditioner  $M_H^{-1}$ , while the second-order scheme (4) is used to build the matrix  $A_H$  for the residual calculation.

$$r_H = \mathcal{I}_h^H r, \tag{9}$$

$$A_H u_H = r_H, \tag{10}$$

$$u_h = \mathcal{I}_H^h u_H, \tag{11}$$

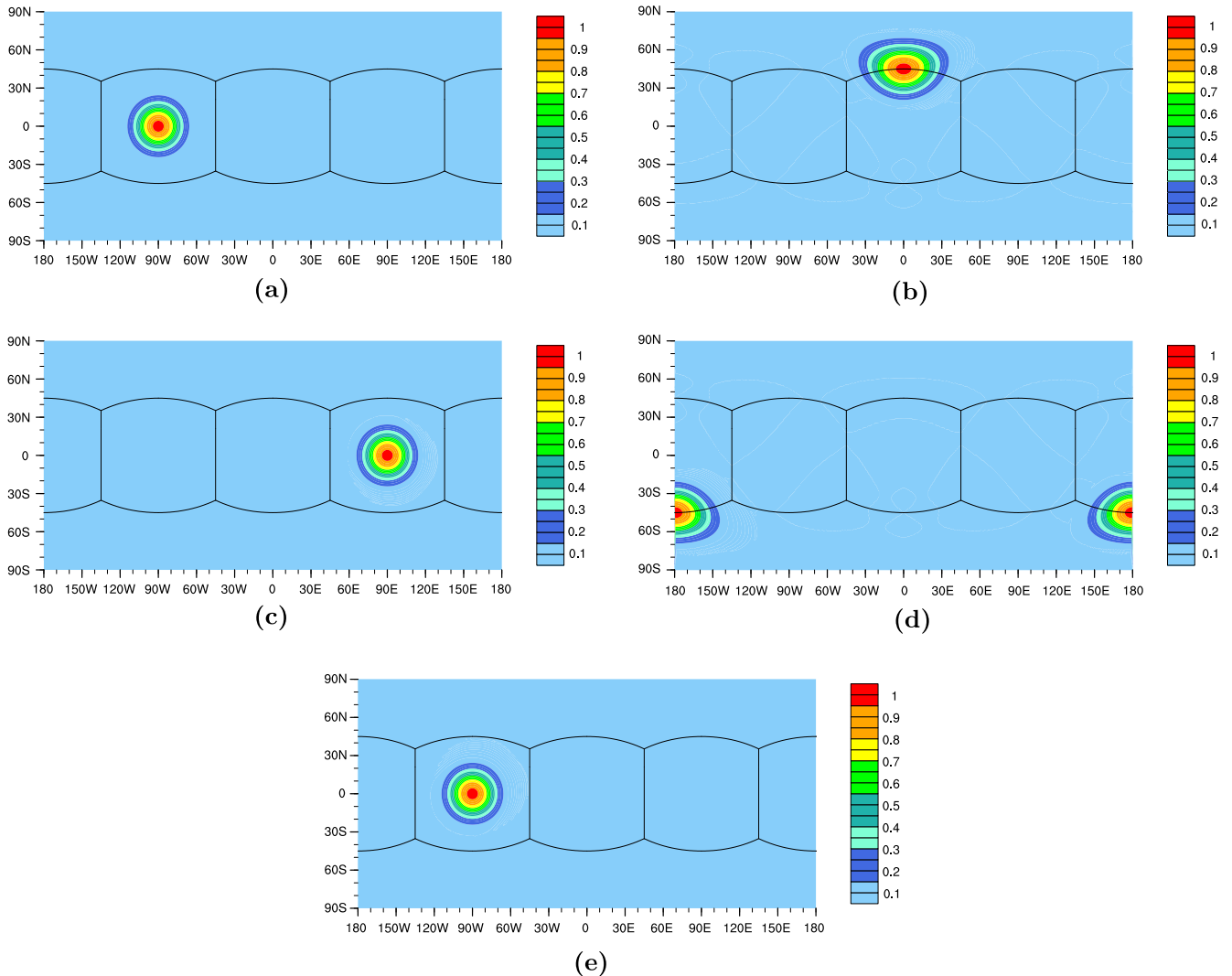


Fig. 2. Contour plots of Case-1. The problem is solved on a  $512 \times 512 \times 6$  mesh. (a) the initial scalar field  $\phi_0$ ; (b) the numerical solution at day three; (c) the numerical solution at day six; (d) the numerical solution at day nine; and (e) the numerical solution at day twelve.

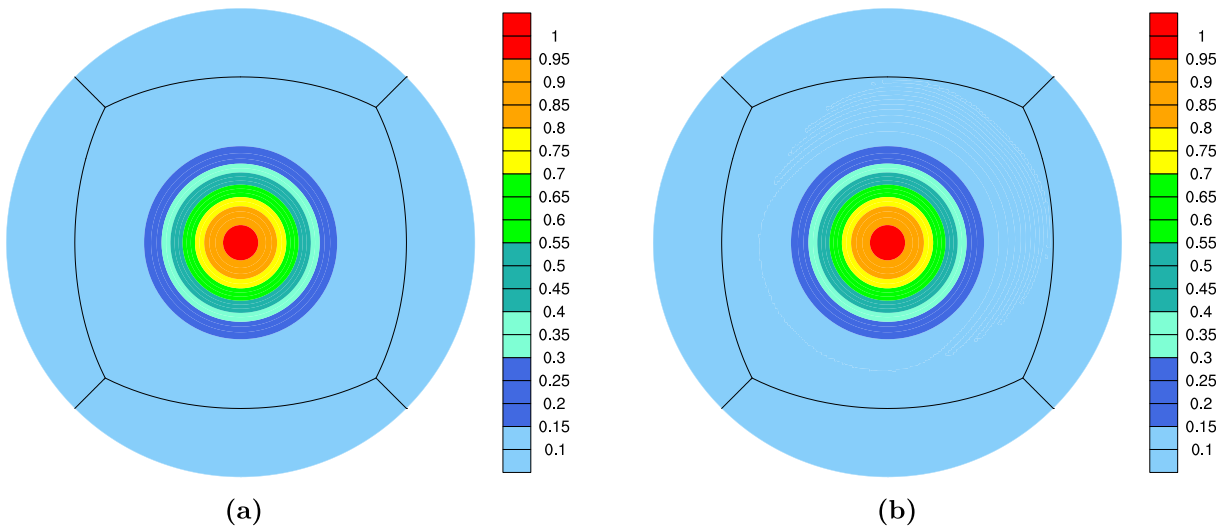


Fig. 3. Contour plots of Case-1 on the sphere. The problem is solved on a  $512 \times 512 \times 6$  mesh. (a) the initial scalar field  $\phi_0$  and (b) the numerical solution at the final time.

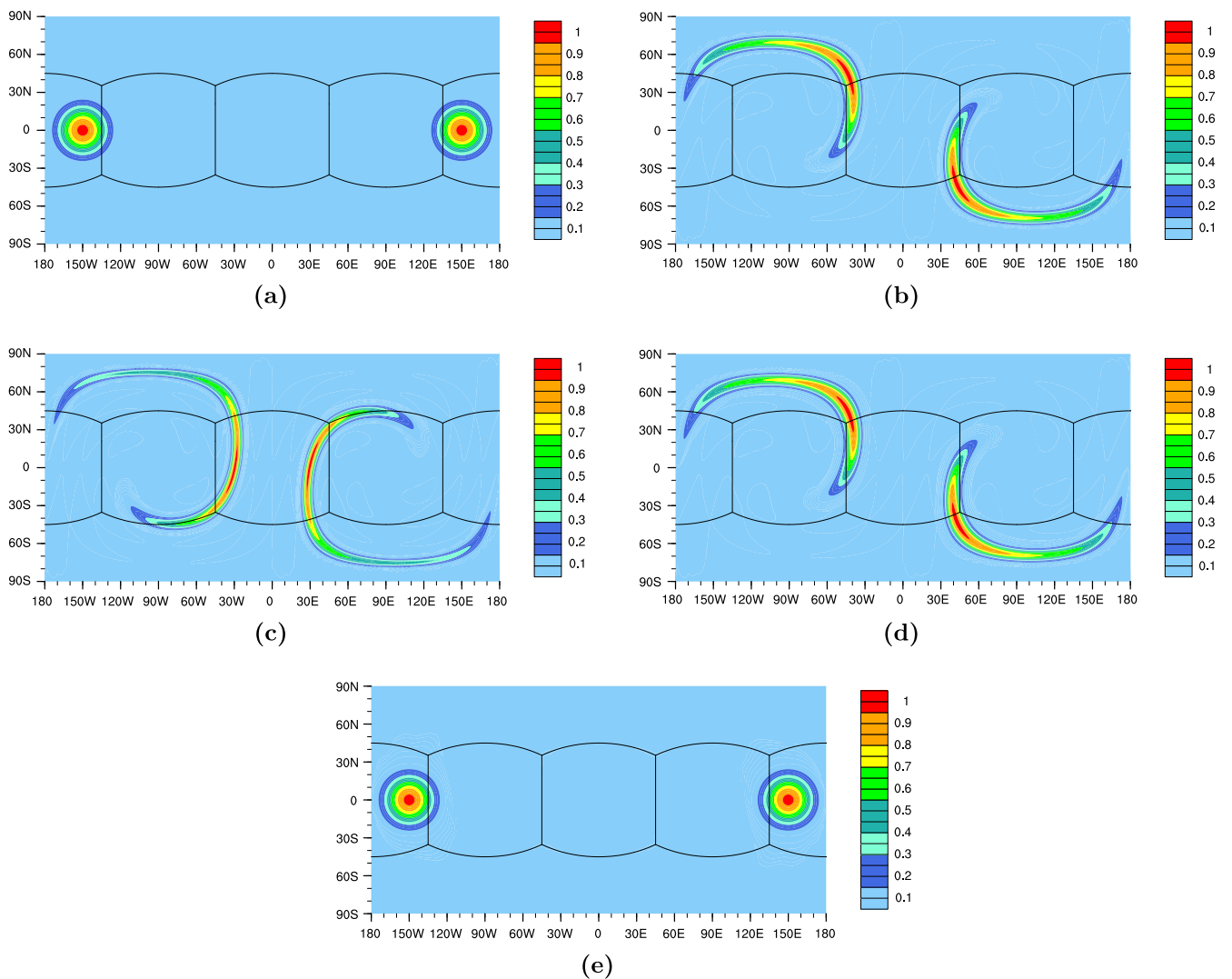


Fig. 4. Contour plots of Case-2. The problem is solved on a  $512 \times 512 \times 6$  mesh. (a) the initial scalar field  $\phi_0$ ; (b) the numerical solution at day three; (c) the numerical solution at day six; (d) the numerical solution at day nine; and (e) the numerical solution at day twelve.

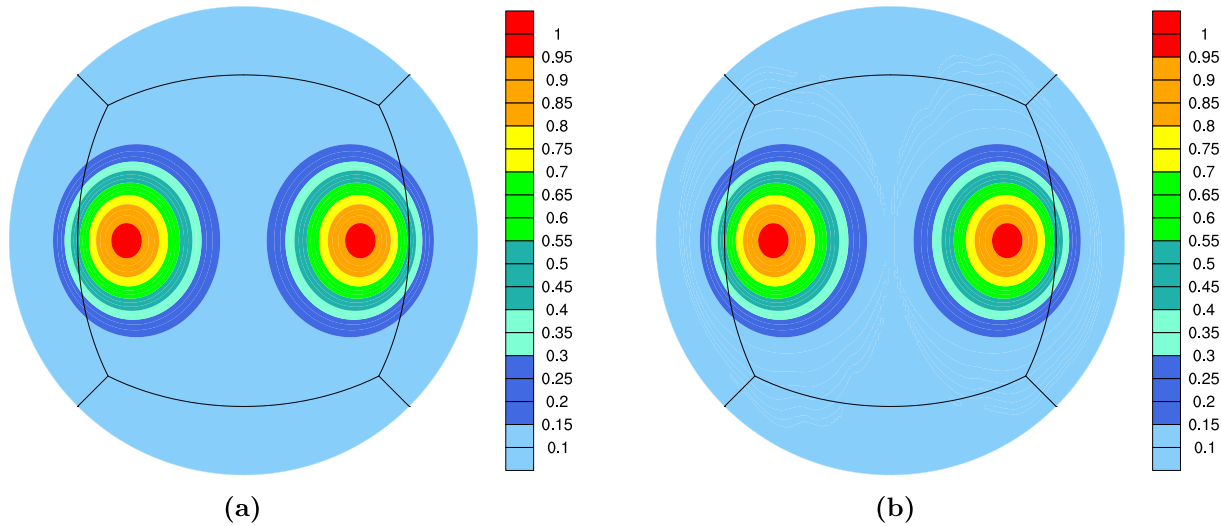


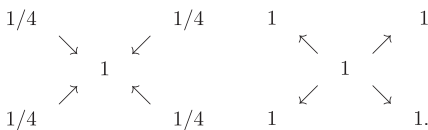
Fig. 5. Contour plots of Case-2 on the sphere. The problem is solved on a  $512 \times 512 \times 6$  mesh. (a) the initial scalar field  $\phi_0$  and (b) the numerical solution at the final time.

Table 1

Effect of overlapping size  $\delta$  for Case-2 with a fixed  $4096 \times 4096 \times 6$  mesh. The time step size is fixed to  $\Delta t = 0.2$  days, the simulation is stopped at day four. The calculation is carried out with 768 processors for the one-level and two-level preconditioners.

$\delta$	One-level		$H = 2h$		$H = 4h$		$H = 8h$	
	GMRES	Time	GMRES	Time	GMRES	Time	GMRES	Time
0	86.3	226.1	29.7	178.1	48.3	172.4	64.0	191.1
1	73.0	220.8	25.6	191.7	42.3	175.3	57.1	205.0
2	72.4	215.8	24.9	192.2	42.5	171.5	57.4	197.3
4	71.5	213.5	24.8	194.6	41.7	180.0	56.2	202.3
6	70.3	219.2	24.6	198.8	41.6	184.1	55.4	213.0
8	70.0	229.2	24.4	205.7	41.4	193.1	54.8	222.9

Another consideration in designing the two-level method is how to formulate the restriction matrix  $\mathcal{I}_h^H$  in (9) and the interpolation matrix  $\mathcal{I}_H^h$  in (11). With the cell-centered finite volume scheme, it happens that a coarse mesh point does not coincide with any fine mesh point; see Fig. 1. We find that the overall method may not converge, if the operators are not chosen properly. In this paper, we build  $\mathcal{I}_h^H$  and  $\mathcal{I}_H^h$  for the case of the coarse-to-fine mesh ratio 1:2, using the following weights



The weights for other coarse-to-fine mesh ratios are defined in a similar way.

Table 2

Effect of time step size  $\Delta t$  for Case-2 with a fixed  $4096 \times 4096 \times 6$  mesh. The simulation is stopped at day four. The calculation is carried out with 768 processors for the one-level and two-level preconditioners.

$\Delta t$	Steps	One-level		$H = 2h$		$H = 4h$		$H = 8h$	
		GMRES	Time	GMRES	Time	GMRES	Time	GMRES	Time
0.025	160	19.4	879.9	10.2	856.1	12.5	745.5	14.5	854.5
0.05	80	34.0	569.6	14.8	544.3	19.5	453.4	24.3	532.2
0.1	40	53.8	392.8	20.6	334.3	31.1	293.6	41.8	343.4
0.5	8	110.3	116.8	38.0	104.6	57.8	87.7	86.5	109.2
1.0	4	199.8	102.9	84.5	89.3	105.0	67.2	153.2	81.6

### 4. Numerical results

The algorithms described in the previous sections are implemented based on the Portable Extensible Toolkit for Scientific computation (PETSc) [1]. All computations are performed on a Dell PowerEdge C6100 supercomputer located at the University of Colorado Boulder. Each node contains 24 GB local memory and two hex-core 2.8 Ghz Intel Westmere processors. The nodes are interconnected via a non-blocking QDR Infiniband high performance network. We focus on the parallel performance of the proposed methods, including the effect on using different parameters in the preconditioner, the robustness of the method when the time step size is large, and the strong and weak scalabilities of the algorithm.

#### 4.1. Test cases

In this subsection, we introduce two published benchmark problems that are used later for numerical tests.

##### 4.1.1. Case-1: Solid-body advection of a cosine bell

The first test case describes a steady and non-divergent wind field that transports a tracer field around a great circle without distortion [6,3,7,17]. The velocity field is set as

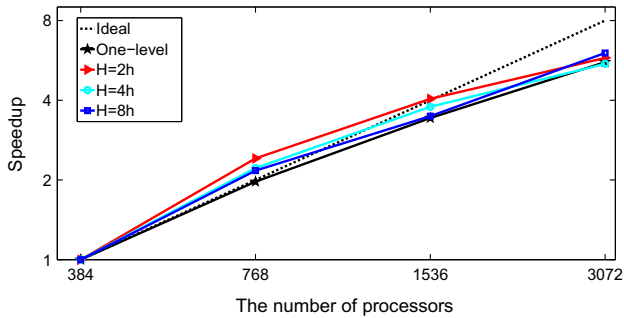
$$\begin{cases} u(\lambda, \theta, t) = u_0(\cos \alpha \cos \theta + \sin \alpha \cos \lambda \sin \theta), \\ v(\lambda, \theta, t) = -u_0 \sin \alpha \sin \lambda, \end{cases}$$

where  $\lambda$  is the longitude,  $\theta$  is the latitude,  $u_0$  is the reference wind velocity, and  $\alpha$  is the angle between the polar axis and the axis of the solid body rotation ( $\alpha = 0$  is for pure zonal flow and  $\alpha = \pi/2$  is for over-the-pole flow). Here, we use  $\alpha = \pi/4$ , which is

**Table 3**

Effect of different number of processors  $N_p$  for **Case-1** and **Case-2** with a fixed  $4096 \times 4096 \times 6$  mesh. The time step size is fixed to  $\Delta t = 0.2$  days, the simulation is stopped at day four.

$N_p$	One-level		$H = 2$ h		$H = 4$ h		$H = 8$ h	
	GMRES	Time	GMRES	Time	GMRES	Time	GMRES	Time
<i>Case-1</i>								
384	134.4	803.9	31.6	585.6	66.0	603.8	102.7	778.3
768	133.6	408.0	32.3	242.8	67.3	272.3	103.5	358.8
1536	134.5	234.5	31.9	144.6	67.6	159.8	106.7	223.4
3072	137.1	144.3	32.4	101.3	67.7	110.0	107.3	129.0
<i>Case-2</i>								
384	75.4	409.7	25.5	344.3	39.8	308.9	54.6	353.2
768	73.0	228.9	25.6	191.7	42.3	175.3	57.1	205.0
1536	79.4	143.7	24.6	109.2	43.4	109.0	58.1	128.7
3072	79.4	93.9	25.7	82.6	45.5	79.5	59.7	88.3



**Fig. 6.** Strong scalability results for **Case-1** with a fixed  $4096 \times 4096 \times 6$  mesh. The time step size is fixed to  $\Delta t = 0.2$  days, the simulation is stopped at day four.

challenging for the test on the cubed-sphere, and  $u_0 = 2\pi/T$  with the final time  $T = 12$  days means that a full rotation around the earth finishes every twelve days. A cosine bell is defined as:

$$\phi_0(\lambda, \theta) = \begin{cases} 0.1 + 0.9h_0(\lambda, \theta) & r_0 < r, \\ 0.1 & \text{otherwise,} \end{cases} \quad (13)$$

where  $r = 1/2$  is the base radius of the bells and

$$h_0(\lambda, \theta) = \frac{1}{2}(1 + \cos(\pi r_0/r)), \quad \text{if } r_0 < r,$$

with  $r_0$  being the great-circle distance between  $(\lambda, \theta)$  and  $(\lambda_0, \theta_0) = (-\pi/2, 0)$ :

$$r_0(\lambda, \theta) = \arccos(\sin \theta_0 \sin \theta + \cos \theta_0 \cos \theta (\lambda - \lambda_0)).$$

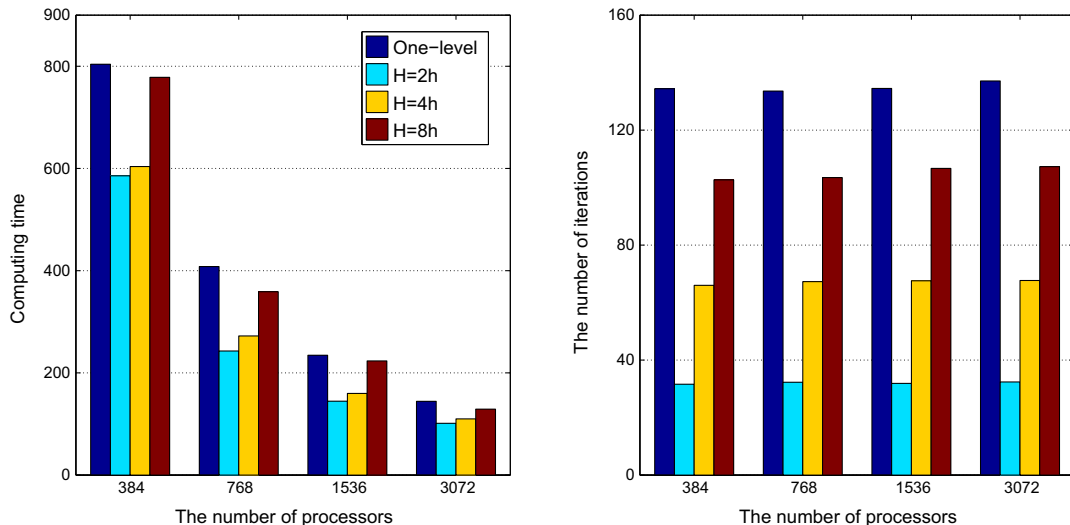
The solid-body rotation of a cosine bell along a great-circle trajectory is a widely used benchmark test, which has a final solution that is identical to the initial condition. Numerical solutions are obtained by using the implicit method, as shown in **Figs. 2 and 3**. We find that the final solution is in good agreement with the initial condition obtained by our method.

**4.1.2. Case-2: Advection of the double cosine bells**

The second test case describes cosine bells with a strongly deformational non-divergent flow [6,8]. The two cosine bells (13) are centered at  $(\lambda_0, \theta_0) = (5\pi/6, 0)$  and  $(-5\pi/6, 0)$ , respectively. The zonal and meridional velocity components in the flow are

$$\begin{cases} u(\lambda, \theta, t) = u_0 \sin^2(\lambda) \sin(2\theta) \cos(\pi t/T), \\ v(\lambda, \theta, t) = u_0 \sin(2\lambda) \cos(\theta) \cos(\pi t/T), \end{cases}$$

respectively, where the final time  $T = 12$  days and  $u_0 = 10/T$ . In the simulation, the two cosine bells follow the complex trajectories of the flow and undergo severe deformation. Moreover, the flow reverses its course at half-time and the scalar fields return to their initial positions and shapes at the final time, which makes the exact solution available at the end of the simulation, as shown in **Figs. 4 and 5**.



**Fig. 7.** Strong scalability results for **Case-1** with different number of processors  $N_p$ . The mesh is  $4096 \times 4096 \times 6$ .

#### 4.2. Performance and parallel scalability

The parallel performance of the domain decomposition preconditioner is investigated. In the one-level method, the second-order discretization is used for the residual calculation and the construction of the RAS preconditioner. In the two-level method, both the second-order and the first-order methods are evaluated using the algorithm introduced in Section 3, the former is used for the residual calculation on both the fine and coarse levels, and the latter is used to build the RAS preconditioner on the coarse level. Moreover, in the two-level preconditioner, the size of the coarse mesh has a strong impact on the efficiency and parallel scaling of the method, using a relatively fine coarse mesh helps reduce the total number of linear iterations, but at the same time the total compute time may increase. An important implementation detail to consider in designing the two-level method is to balance the effects of preconditioning and the computing time of the coarse solve. Let  $h$  be the size of the fine mesh and  $H$  the size of the coarse mesh. In the experiments, we compare the following algorithms using different coarse-to-fine mesh ratios:

- *One-level*: one-level additive Schwarz method (7) is used as the preconditioner.
- $H = 2h$ : two-level multiplicative Schwarz method (8) is used as the preconditioner with the coarse-to-fine mesh ratio 1:2.
- $H = 4h$ : two-level multiplicative Schwarz method (8) is used as the preconditioner with the coarse-to-fine mesh ratio 1:4.
- $H = 8h$ : two-level multiplicative Schwarz method (8) is used as the preconditioner with the coarse-to-fine mesh ratio 1:8.

In the tests, GMRES(30) and FGMRES(30) are used to solve the linear systems on the coarse and the fine meshes, respectively. We use  $10^{-5}$  as the absolute stopping condition for all linear solves, except for the coarse solve of the two-level preconditioner, for which we use  $10^{-1}$  as the relative stopping tolerance. The subdomain problems are solved with a sparse ILU factorization with four levels of fill-ins.

Throughout this subsection, “GMRES” denotes the average number of iterations per time step, and “Time” denotes the total computing time in seconds.

##### 4.2.1. Impact of the overlapping parameter

For the Schwarz preconditioner, the overlapping size  $\delta$  is an important parameter that influences the effect of the preconditioner. From the theory of the overlapping Schwarz method, larger

overlap often implies a faster convergence in terms of GMRES iterations. However, larger overlap also means larger subdomain problems and more information transfers between subdomains; as a result, the overall compute time may increase. Table 1 shows the effect of the overlapping parameter for **Case-2**. We observe from the table that the methods converge better as the overlap increases and on the other hand a larger overlap may also increase inter-process communications.

##### 4.2.2. Impact of time step sizes

Next, to show the unconditional stability of the fully implicit scheme, we solve the same problem with gradually increased time step size  $\Delta t$ . Table 2 shows the number of iterations and the computing time of the fully implicit method with respect to different time step sizes for **Case-2**. As  $\Delta t$  increases, the total number of GMRES iterations increases, while the computing time decreases due to the reduced number of time steps. Since we use a fully implicit method in which the time step size is no longer constrained by the CFL condition, the results clearly indicate that the one-level and two-level methods are convergent for even very large value of  $\Delta t$ .

##### 4.2.3. Impact of the number of processors

Besides the robustness with respect to the overlapping parameter  $\delta$  and time step sizes  $\Delta t$ , scalability is another important concern in the design of solution algorithms for the tracer transport problem. As mentioned in [15,16], the domain decomposition method is still not well understood for hyperbolic problems such as the tracer transport problem, due to the lack of ellipticity. Therefore we numerically investigate the strong and weak scalabilities of the proposed algorithms for **Case-1** and **Case-2** in terms of the computing time and GMRES iterations. We first study the strong scalability of the methods by using a fixed mesh  $4096 \times 4096 \times 6$  and also a fixed time step size  $\Delta t = 0.2$  days in Table 3. We see that the algorithms are scalable with respect to the number of processors, the number of GMRES iterations varies mildly with different  $N_p$ . In Fig. 6, we report the average computing time and the average number of GMRES iterations per time step with respect to the number of processors. It is important to note that the timing results obtained by using the two-level preconditioners are always better than the results obtained with the one-level preconditioner (see Fig. 7).

We further test our algorithms in terms of the weak scalability with a fixed  $192 \times 192$  mesh per processor. We refine the mesh and increase the number of processors simultaneously to keep

**Table 4**  
Weak scalability results for **Case-1** and **Case-2** with a fixed  $192 \times 192$  mesh per processor. The time step size is fixed to  $\Delta t = 0.2$  days, the simulation is stopped at day four. The total degree of freedom for the largest simulation is  $4224 \times 4224 \times 6 = 107,053,056$ .

$N_p$	Mesh	One-level		$H = 2h$		$H = 4h$		$H = 8h$		
		GMRES	Time	GMRES	Time	GMRES	Time	GMRES	Time	
<i>Case-1</i>										
24	$384 \times 384 \times 6$	10.9	16.9	6.8	15.3	8.7	14.9	9.7	15.3	
96	$768 \times 768 \times 6$	18.1	20.6	9.7	18.4	13.4	21.2	16.0	22.9	
384	$1536 \times 1536 \times 6$	37.3	45.0	16.8	36.5	24.2	38.0	29.8	38.9	
864	$2304 \times 2304 \times 6$	58.8	68.9	22.9	56.1	35.7	56.2	45.3	62.6	
1536	$3072 \times 3072 \times 6$	89.1	95.4	27.2	73.5	47.9	75.9	67.7	92.3	
1944	$3456 \times 3456 \times 6$	109.4	118.5	29.2	85.3	54.3	87.6	79.8	110.7	
2904	$4224 \times 4224 \times 6$	146.7	163.5	32.8	106.1	68.6	115.0	115.1	135.0	
<i>Case-2</i>										
24	$384 \times 384 \times 6$	9.9	13.4	7.5	16.2	8.4	14.6	8.6	14.4	
96	$768 \times 768 \times 6$	16.0	16.7	10.2	19.1	12.5	17.8	14.0	18.5	
384	$1536 \times 1536 \times 6$	25.6	26.5	14.5	26.9	18.6	24.9	22.1	26.8	
864	$2304 \times 2304 \times 6$	37.6	41.5	17.8	40.1	25.7	37.6	31.3	40.5	
1536	$3072 \times 3072 \times 6$	53.5	65.9	20.9	53.3	33.5	50.6	44.0	60.6	
1944	$3456 \times 3456 \times 6$	64.7	89.5	22.6	58.3	38.0	53.1	50.8	62.8	
2904	$4224 \times 4224 \times 6$	89.3	111.4	25.7	86.8	46.9	80.8	64.9	90.1	

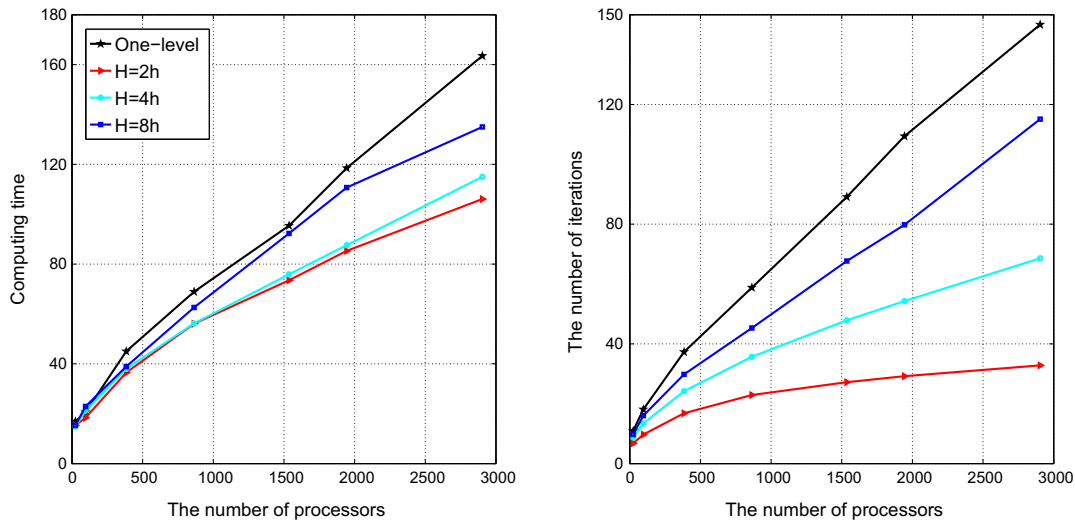


Fig. 8. Weak scalability results for **Case-1** with a fixed  $192 \times 192$  mesh per processor. The time step size is fixed to  $\Delta t = 0.2$  days, the simulation is stopped at day four.

the number of unknowns per processor as a constant. The test starts with a  $384 \times 384 \times 6$  mesh for  $N_p = 24$  and ends up with a  $4224 \times 4224 \times 6$  mesh for  $N_p = 2904$ , as shown in Table 4. For the fully implicit methods, the condition number of the preconditioned matrix becomes larger when we refine the mesh and increase  $N_p$ , leading to the growth of the number of GMRES iterations, especially for the one-level method. If we focus on **Case-1**, observing from Fig. 8, we remark that the number of linear iterations for the case of  $H = 2h$  is much smaller than for other cases, and the total solution time behaves similarly. Meanwhile, for **Case-2**, we find that the two-level method with  $H = 4h$  shows the best performance in terms of the computing time.

## 5. Conclusion

In this paper, we developed a parallel two-level domain decomposition method with mixed order discretizations for the tracer transport problem on the cubed-sphere mesh, based on a fully implicit finite volume discretization. Since the scalability of the proposed method depends heavily on the design of the preconditioner, we used the idea of mixed order discretizations to build a class of two-level restricted Schwarz methods. In the approach, we built the Schwarz preconditioner in a multiplicative manner, in which the preconditioner on the coarse mesh is based on the first-order spatial discretization while the second-order scheme is used to construct the preconditioner on the fine mesh. Finally, we tested the algorithm for two benchmark problems to show the robustness and accuracy on high resolution meshes with more than one hundred million unknowns. The numerical experiments indicate that the method performs well in terms of both strong and weak scalabilities with up to 3072 processors.

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