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TURBULENT PARAMETERIZATION FOR CCATT-BRAMS BY GPU

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ABSTRACT

A strategy to increase the computer code performance is to employ of hybrid computing, combining CPU with GPU or/and FPGA. Two turbulent parameterizations, Smagorinsky and Mellor-Yamada, were codified on GPU for the CCATT-BRAMS code, using the CUDA standard. The results show a good speed-up for the model.

RESUMO

Uma estratégia para aumentar o desempenho de códigos computacionais é o uso de computação híbrida, que combina CPU com GPU e/ou FPGA. Duas parametrizações de turbulência, Smagorinsky e Mellor-Yamada, foram codificadas em GPU para o código CATT-BRAMS:. O padrão CUDA foi usado na implementação. Os resultados mostram um aumento significativo no speed-up do modelo.

1. INTRODUCTION

The environmental prediction model CCATT-BRAMS (Freitas et al., 2009; 2013) is a complex computer code, and it is operationally employed by the CPTEC-INPE. This model allows atmospheric simulation with emission, transport, pollutant dispersion, and chemical reactions. The CCATT-BRAMS is already prepared to run on parallel machines. But, depending on the model resolution, this is a very expensive code. This motivates the research to improve its performance. Two parameterizations in the CCATT-BRMAS are codified on GPU: Smagorinsky (1963), and Mellor-Yamada (1982).

1.1 Smagorinsky's parameterization

The turbulent fluxes are parameterized using the theory of gradient-flux. The Reynolds tensors are given as following:

$$\overline{U'_i U'_j} = -(K_m)_{ij}(D)_j \tag{1a}$$

$$(D)_{j} = \partial \overline{u}_{j} / \partial x_{j} + \partial \overline{u}_{j} / \partial x_{j}$$

$$(1b)$$

$$K_m = (CS_2\Delta z)^2 \left[|D_v| + H(N) \right] f(R)$$
(1c)

where cs_z is a fitting coefficient, Δz is the discretization for vertical grid, and $|D_v|$ is the magnitude of the deformation tensor for vertical direction. H(N) and $f(R_i)$ are given by

$$H(N) = \sqrt{\max(0, -N^2)}$$
 (2a)

$$f(R_i) = \max\{0 ; 1 - (K_h/K_m)R_i\}$$
 (2b)



with N being the Brunt-Väisälä frequency, and (K_h/K_m) is the ratio between the heat and momentum eddy diffusivities.

1.2 Mellor-Yamada's parameterization

This is a 2.5 closure scheme for turbulence, where the third order tensors are parameterized, with a prognostic equation for the turbulent kinetic energy (e):

$$\frac{\partial \overline{\psi_{i}\phi_{j}'}}{\partial t} = f\left(\overline{\psi_{i}\phi_{j}'\zeta_{k}'}, \overline{\psi_{i}\phi_{j}'}, \psi_{m}, t\right)$$
(3a)

$$\overline{\psi_i \phi_j' \zeta_k'} \approx K_p \nabla \overline{\psi_i \phi_j'}$$
(3b)

where ψ , ϕ , $\zeta = u$, v, w, q, θ (wind components, moisture, and temperature). Key parameters are a length scale l, and TKE (e):

$$K_{p} = S_{p}I\sqrt{2e} , \quad I = \frac{\kappa(\mathbf{z} + \mathbf{z}_{0})}{1 + \kappa(\mathbf{z} + \mathbf{z}_{0})I_{\infty}}, \quad I_{\infty} = 0.1 \frac{\int_{0}^{d} \mathbf{z}\sqrt{2e} d\mathbf{z}}{\int_{0}^{d} \sqrt{2e} d\mathbf{z}}$$

$$(4)$$

where S_P are constants suggests by Mellor-Yamada (1982).

2. CUDA PROGRAMMING ON GPU

CUDA (Compute Unified Device Architecture) (Nvidia: CUDA, 2012) was developed for the parallel processing for the GPUs produced Nvidia Corporation. The CUDA programming could be understood as an extension of languages like C, C++, and Fortran, adding qualifiers to functions and data, producing kernels for execution. The job on a kernel is to be divided among thousands of threads, organized into blocks, and grids, with dimensions *blockDim* and *gridDim*. The kernel uses indexes *blockIdx* and *threadIdx* for defining the job to a thread.

The turbulence routines were codified on Nvidia GPUs: Fermi GTX-580 (512 cores), and Kepler GTX-680 (1536 cores).

3. RESULTS AND FINAL COMMENTS

Two horizontal resolutions were evaluated: 40 km (grid points number: 95×48×38), and 20 km (194×100×40). Figure 1 shows the speed-up results to Smagorinsky's approach for both resolutions, and for Fermi and Kepler architectures.

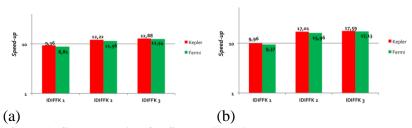


Figure 1: GPU execution for Smagorinsky's approach under two resolutions: (a) 40 km, (b) 20 km.

Similarly, the speed-up for the Mellor-Yamada scheme is presented in Figure 2, also considering two resolutions, and Fermi and Kepler architectures performance.

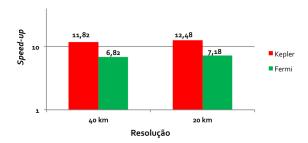


Figure 2: GPU execution for Mellor-Yamada's approach under two resolutions: 40 km, and 20 km.

For the lower resolution worked here, similar speed-up results were obtained for both GPU architectures – the gain for the Mellor-Yamada approach has superior speed-up than the Smagorinsky's parameterization. The results show a tendency to improve the speed-up for a finer resolution.

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