Introduction to High Performance Scientific Computing

Autumn, 2016

Lecture 14

Vectorizing code

In general (Python, Fortran, Matlab,...), avoid for loops and *vectorize* calculations involving arrays.

Example:

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Example:

- Vectorized code will usually be faster, sometimes much faster
- Exception: parallelizing Fortran code with OpenMp: vectorized code → loops → parallel loops

Today

Programming example: from PDE → algorithm → serial code → parallel code

Task: Compute temperature distribution in a room

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Governing equation: Heat equation (diffusion equation):

$$\frac{\partial T}{\partial t} = \alpha \nabla^2 T + S(\mathbf{x},t)$$

$$T(\mathbf{x},t=0) = f(\mathbf{x})$$
 Initial condition

Here, S is a heat source. Boundary conditions should also be specified as appropriate.

Problem: given the source, initial condition, and boundary conditions, solve for the temperature distribution, $T(\mathbf{x},t)$

Today: 1-D problem

$$\frac{\partial T}{\partial t} = \frac{\partial^2 T}{\partial x^2} + S(x, t)$$

$$T(x, t = 0) = f(x)$$

$$T(x = 0, t) = a(t), \ T(x = 1, t) = b(t)$$

$$0 \le x \le 1$$

Initial condition

Boundary conditions

Today: 1-D problem

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 Initial condition
$$T(x=0,t) = a(t), \ T(x=1,t) = b(t)$$
 Boundary conditions
$$0 < x < 1$$

First consider steady problem, e.g., S = S(x), a and b are constants:

$$\frac{\partial^2 T}{\partial x^2} + S(x,t) = 0 \qquad \text{Poisson equation}$$

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Notes:

- 1. This is an extremely simple problem, easy to write down the analytical solution
- 2. No need to use compiled language
- 3. Certainly no need to parallelize
- 4. But what about two-dimensional or three-dimensional problems?
 - Then, the picture changes considerably!
- 5. We are just considering the 1-D problem for illustrative purposes

First consider steady problem, e.g., S = S(x), a and b are constants:

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 Poisson equation

Numerical method:

1. Discretize the derivative:

$$\frac{\partial^2 T}{\partial x^2} \approx \frac{T_{i+1} - 2T_i + T_{i-1}}{\Delta x^2}$$

$$x_i = i * \Delta x, \ i = 1, 2, ..., N$$

$$(N+1) * \Delta x = 1$$

2nd-order, centered scheme

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2nd-order, centered scheme

$$x_i = i * \Delta x, \ i = 0, 1, 2, ..., N + 1$$

 $(N+1) * \Delta x = 1$

With boundary conditions: $T_0 = T_a, \ T_N = T_b$

Equation for T_i:
$$\frac{T_{i+1}-2T_i+T_{i-1}}{\Delta x^2}=-S_i$$

In matrix form: AT = b

$$A = \begin{bmatrix} -2 & 1 & 0 & 0 & 0 & \dots & 0 \\ 1 & -2 & 1 & 0 & 0 & \dots & 0 \\ 0 & 1 & -2 & 1 & 0 & \dots & 0 \\ \vdots & \ddots & \ddots & \ddots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 1 & -2 & 1 \\ 0 & \dots & 0 & 0 & 0 & 1 & -2 \end{bmatrix}, b = \frac{1}{\Delta x^2} \begin{bmatrix} -\Delta x^2 T_a - S_1 \\ -S_2 \\ \vdots \\ -S_i \\ \vdots \\ -S_{N-1} \\ -\Delta x^2 T_b - S_N \end{bmatrix}$$

- In 1-D, this is just a tridiagonal system of equations
- Easy to solve directly (with, say, DGTSV)

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- Then, direct solution becomes very expensive for large N
- Iterative methods are a popular alternative

- In two or three dimensions, A loses it simple banded structure
- Then, direct solution becomes very expensive for large N
- Iterative methods are a popular alternative
- Basic idea: rewrite Ax=b as $A_1x = A_2x + b$
- Choose A₁ so that it is easy to invert, then solve iterative system:
- $A_1 x^{k+1} = A_2 x^k + b$
 - Requires guess, x⁰

Jacobi iteration

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- Choose A₁ so that it is easy to invert, then solve iterative system:
- $A_1 x^{k+1} = A_2 x^k + b$
 - Requires guess, x⁰
- Jacobi iteration: Choose A₁ to be diagonal matrix (main diagonal of A):

$$\frac{T_{i+1}^{k-1} - 2T_i^k + T_{i-1}^{k-1}}{\Delta x^2} = -S_i$$

$$T_i^k = \frac{\Delta x^2}{2} S_i + \frac{1}{2} \left(T_{i+1}^{k-1} + T_{i-1}^{k-1} \right)$$

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Main algorithm, easy to code!

Jacobi iteration in Fortran

- Plan:
 - Set parameters: a, b, n, tol
 - Construct grid x_i
 - Construct source function, S(x), initialize T=T(x,t=0)
 - Iterate using formula below
 - Each iteration check if |Tk-Tk-1| < tol

$$T_i^k = \frac{\Delta x^2}{2} S_i + \frac{1}{2} \left(T_{i+1}^{k-1} + T_{i-1}^{k-1} \right)$$

Jacobi iteration in Fortran

One Fortran trick: set variables to be dimension(0:N+1)

•
$$x(0)=0$$
, $x(N+1)=1$, $T(0)=a$, $T(N+1)=b$

• Then, easy to compute T₁ using:

$$T_i^k = \frac{\Delta x^2}{2} S_i + \frac{1}{2} \left(T_{i+1}^{k-1} + T_{i-1}^{k-1} \right)$$

Jacobi iteration in Fortran

- One Fortran trick: set variables to be dimension(0:N+1)
 - x(0)=0, x(N+1)=1, T(0)=a, T(N+1)=b
 - Then, easy to compute T₁ using:

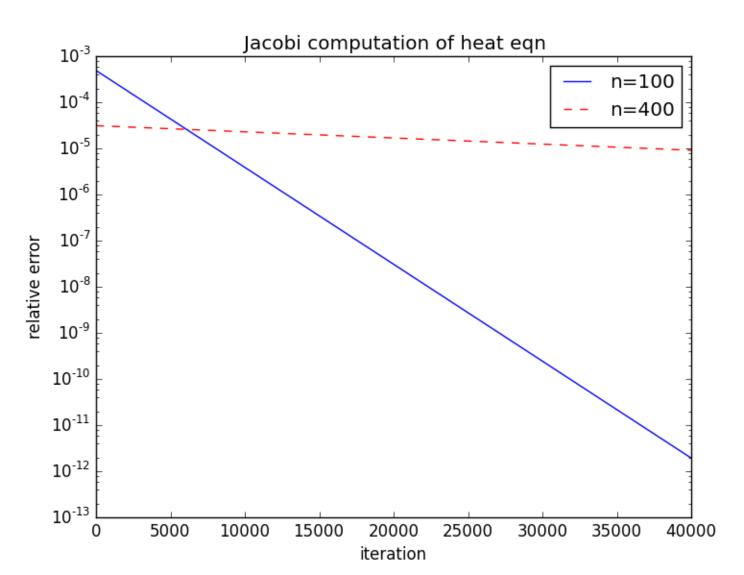
$$T_i^k = \frac{\Delta x^2}{2} S_i + \frac{1}{2} \left(T_{i+1}^{k-1} + T_{i-1}^{k-1} \right)$$

Core part of code (see jacobi1s.f90):

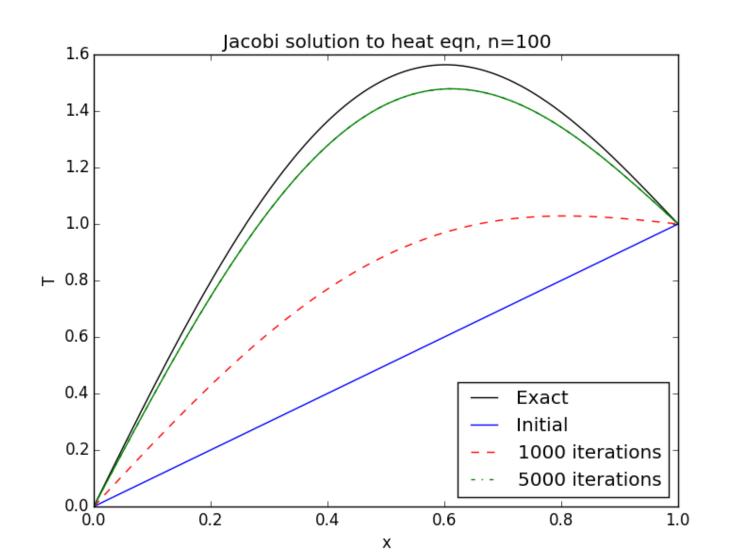
```
do k1=1,kmax
```

1. Does the solution converge?

1. Does the solution converge? Yes, but very slowly for large n



1. Does it converge to the correct solution? Yes, can check that error $\sim \Delta x^2$



- Jacobi is simplest, but most inefficient iterative solver
- Good illustration of basic ideas
- Better methods: Gauss-Seidel, SOR, conjugate gradient, multigrid

Let's now parallelize the solver with OpenMP

- Look for loops that can be parallelized
- Look for vectorized operations that can be converted to loops that can be parallelized

Serial:

```
Tnew(1:n) = S(1:n)*dx2f + 0.5d0*(T(0:n-1) + T(2:n+1)) !Jacobi deltaT(k1) = maxval(abs(Tnew(1:n)-T(1:n))) !compute relative error
```

Let's now parallelize the solver with OpenMP

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Parallel:

```
dmax=0.d0
!$omp parallel do reduction(max:dmax)
do i1=1,n
        Tnew(i1) = S(i1)*dx2f + 0.5d0*(T(i1-1) + T(i1+1))
        dmax = max(dmax,abs(Tnew(i1)-T(i1)))
end do
!$omp end parallel do
deltaT(k1) = dmax
```

Let's now parallelize the solver with OpenMP

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Serial:

```
do i1=0,n+1
     x(i1) = i1*dx
end do
!-----
!set initial condition
T = (b-a)*x + a
!set source function
S = S0*sin(pi*x)
```

Let's now parallelize the solver with OpenMP

- Look for loops that can be parallelized
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Parallel:

```
!$omp parallel do
do i1=0,n+1
    x(i1) = i1*dx
    T(i1) = (b-a)*x(i1) + a !set initial condition
    S(i1) = S0*sin(pi*x(i1)) !set source function
end do
!$omp end parallel do
```

Parallel Jacobi notes

- Will only see speedup with n > ~20000 (commonly seen in 2D problems)
- See jacobi1s_omp.f90, jacobi1_omp.py
- f2py and OpenMP: f2py --f90flags='-fopenmp' -lgomp -c jacobi1s_omp.f90 -m j1

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$$0 < x < 1$$

Simple (inefficient) approach: method of lines

- 1. Discretize spatial variable → N+2 points beteween 0 and 1
- 2. Solve resulting N ODEs with solver of choice (odeint, ode15s ,...)

Again, we discretize the derivative:

$$\frac{\partial^2 T}{\partial x^2} \approx \frac{T_{i+1} - 2T_i + T_{i-1}}{\Delta x^2}$$

$$x_i = i * \Delta x, \ i = 0, 1, 2, ..., N + 1$$

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Again, we discretize the derivative as:

$$\frac{\partial^2 T}{\partial x^2} \approx \frac{T_{i+1} - 2T_i + T_{i-1}}{\Delta x^2}$$

$$x_i = i * \Delta x, \ i = 0, 1, 2, ..., N + 1$$

$$(N+1) * \Delta x = 1$$

So, we have N ODEs:

$$\frac{dT_i}{dt} = S_i(t) + \frac{T_{i+1} - 2T_i + T_{i-1}}{\Delta x^2}, \ i = 1, 2, ..., N$$

with the boundary conditions substituted in the RHS when needed.

Solving single ODE in python (lecture 6)

Use odeint from scipy.integrate module to solve:

$$\frac{dy}{dt} = -ay$$

- Basic idea: discretize time, t = 0, dt, ..., N*dt, and starting from y(0) march forward in time and compute y(dt), ... y(N*dt)
- odeint chooses the stepsize, dt, so that error tolerances are satisfied
- Need to specify:
 - Initial condition
 - Timespan for integration
 - A Python function which provides RHS of the ODE to odeint
- Look at ode_example.py and lab 4

Solving N ODEs:

$$\frac{dT_i}{dt} = S_i(t) + \frac{T_{i+1} - 2T_i + T_{i-1}}{\Delta x^2}, \ i = 1, 2, ..., N$$

- Will need to provide N initial conditions when calling odeint.
- The python function which provides RHS to odeint will:
 - Take t and $T_1, ..., T_N$ and any other needed parameters as input
 - Return N values for dT/dt as output
- No need for Fortran for 1D problems, but may be faster for two and three dimensions.