

Parallel Algorithms for the Solution and Estimation of DSGE Models

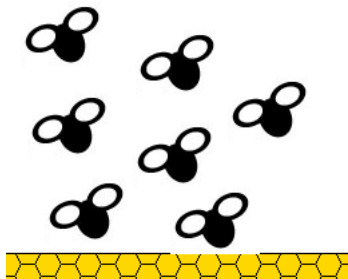
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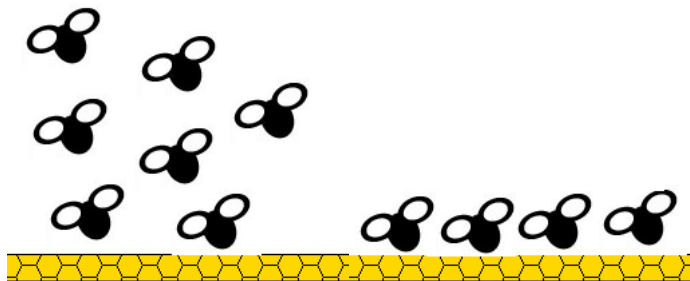
June 27, 2008

Threads Ideal



A thread is a call sequence the executes independently of other threads. It may share underlying system resources with other threads.

Threads Synchronized



Model Class

$$f(s, x, Eh(s, x, e', s', x')) = 0$$

$$s' = g(s, x, e')$$

Specified Functions

f	equilibrium conditions
h	expectation functions
g	state transitions

Variables

s	state variables
x	policy variables
z	expectation variables
e	stochastic shocks

Standard RBC Model for N Countries

$$\max U \{ \{ c_{n,t}, l_{n,t}, i_{n,t} \}_{n=1}^N \}_{t=0}^{\infty}$$

$$U = E_0 \sum_{t=0}^{\infty} \sum_{n=1}^N \beta^t U_{n,t}$$

$$U_{n,t} = \frac{(c_{n,t}^{\theta_n} (1 - l_{n,t})^{1-\theta_n})^{1-\tau_n}}{1 - \tau_n}$$

$$\sum_{n=1}^N (y_{n,t} - c_{n,t} - i_{n,t}) = 0$$

$$y_{n,t} = e^{a_{n,t}} k_{n,t}^{\alpha_n} l_{n,t}^{1-\alpha_n}$$

$$k_{n,t+1} = (1 - \delta_n) k_{n,t} + i_{n,t} - 0.5 \kappa_n i_{n,t}^2$$

$$a_{n,t+1} = \rho_n a_{n,t} + e_{n,t+1}$$

Solution

Solutions can be characterized as policy functions

$$x_t = x^*(s_t)$$

with $x_t = (c_t, l_t)'$ and $s_t = (k_t, a_t)'$

In general no analytic solution is possible. Global polynomial approximation (Judd, 1992) may be used (or linearization, perturbation...).

Approximation Strategy

Policy functions by Chebychev polynomials:

$$x = \Psi(s) = \sum_{i=1}^n c_i \psi_i(s)$$

Expectations over e' via Gauss Quadrature:

$$\begin{aligned} E h(s, x, e', s', x') &= \\ &= \sum_j \omega_j h(s, x^*(s), e'_j, s'_j, x^*(s'_j)) \end{aligned}$$

over optimal nodes and weights. We use the Smolyak operator (Smolyak, 1963) to cope with the curse of dimensionality.

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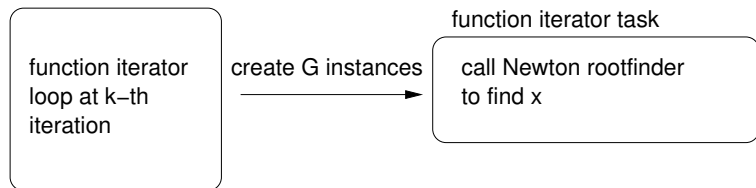
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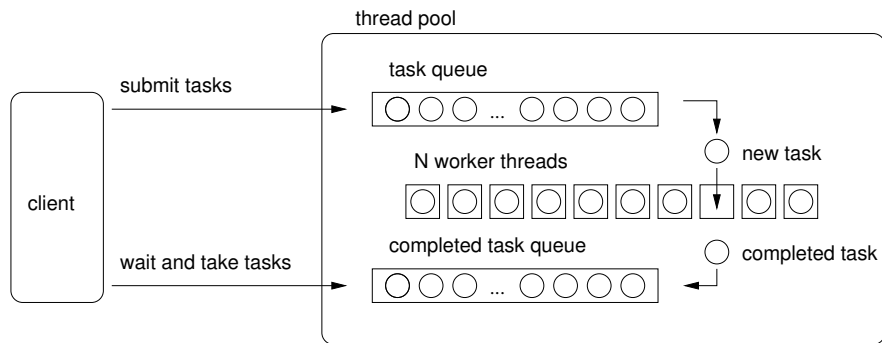
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4. $k=k+1$, go to (1) until $R \approx 0$

Parallelization



Thread Pool Pattern



implemented in *java.util.concurrent* package and many others

State Space

The model solving policy function $x(s)$ can be used to substitute out the policy variables in the state transition equation.

Augmented with a measurement equation we obtain the model's empirical implication for the observables in a nonlinear state space model

$$\begin{aligned} s_t &= g(s_{t-1}, x(s_{t-1}), e_t) = g^*(s_{t-1}, e_t) \\ y_t &= m(s_t) + \epsilon_t. \end{aligned}$$

With distributional assumptions for the state and measurement shocks, e_t and ϵ_t , the equations can be cast in terms of state and measurement densities

$$p(s_t | s_{t-1})$$

$$p(y_t | s_t),$$

respectively.

The filtering approach evaluates the likelihood of this model by the prediction and filtering step:

$$\mathcal{L}(\theta; y_{1:T}) = p(y_{1:T} | \theta) = \prod_{t=1}^T p(y_t | y_{1:t-1}, \theta) = \prod_{t=1}^T l_t.$$

Parallel Metropolis-Hastings Algorithm with Mixing

1. Choose start values $\hat{\theta}_{1,m}$ for all $m = 1, \dots, M$ and b , γ^{GE} for an acceptance ratio of $\approx 30\%$

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2.2.1 Acceptance:

$$\hat{\theta}_{n,m} = \begin{cases} \hat{\theta}_m^* & \text{if } U(0, 1) \leq \frac{\rho(y_{1:t}|\hat{\theta}_m^*)\rho(\hat{\theta}_m^*)}{\rho(y_{1:t}|\hat{\theta}_m)\rho(\hat{\theta}_m)} \\ \hat{\theta}_{n,m} & \text{otherwise} \end{cases}$$

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3. Disregard burn-in draws $\hat{\theta}_{1:J,1:M}$

Simulation Setup for Nonlinear Solution

Cluster with 16 Xeon CPUs at 2.7GHz (but only 14 can be free);
GNU/Linux openSUSE 10.2;
Java virtual machine 1.6.0_06

▶ parameters:

$$\alpha_i = 0.4, \beta = 0.99, \delta_i = 0.02, \kappa_i = 0.01, \rho_i = 0.95, \tau_i = 2$$
$$\theta_i = 0.357, \sigma_{a_i} = 0.007$$

▶ states grid bounds:

$$a_i \in [-0.06; 0.06], \quad k_i \in [15; 30]$$

▶ approximation level = 3, integration level = 3

▶ tolerance for function iterator: 1×10^{-4}

▶ starting values for steady state:

$$a_i = 0, i_i = 0.5, k_i = 23, l_i = 0.3$$

▶ strong convergence tolerance for Newton rootfinder: 1×10^{-10}

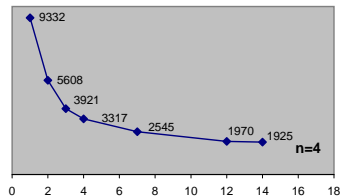
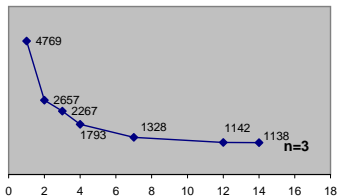
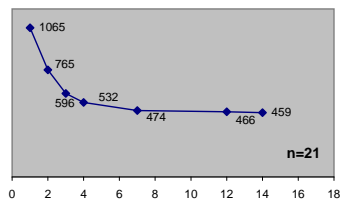
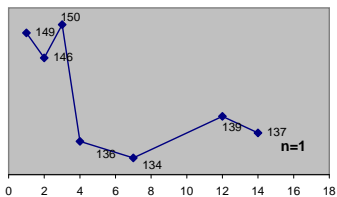
▶ set thread pool size to 1, 2, 3, 4, 7, 12, 14

Amdahl's Law

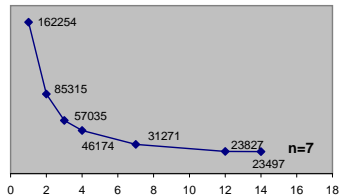
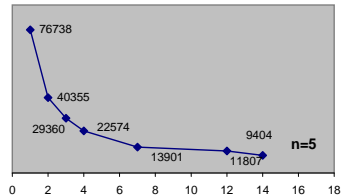
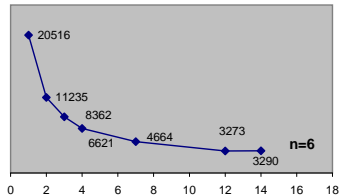
$$\text{sup speedup} = \frac{1}{F + (1 - F)/N}$$

- ▶ F is percentage of system that cannot be parallelized
- ▶ N is number of processors
- ▶ provides a theoretical upper bound for the expected speedup of using N processors instead of one

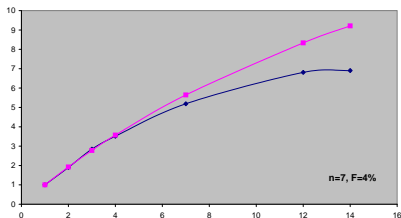
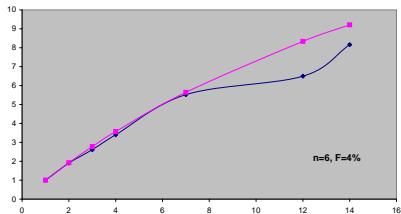
Nonlin. Sol.: Computing Times vs. # Processors



Nonlin. Sol.: Computing Times vs. # Processors cont'd



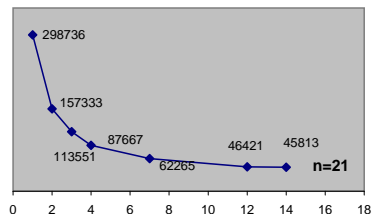
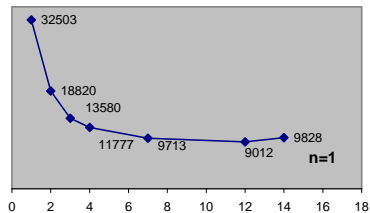
Nonlin. Sol.: Speedup vs. Amdahl's Law



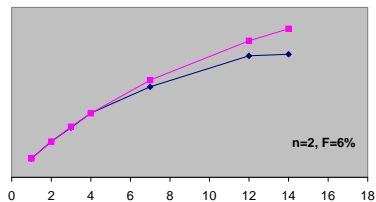
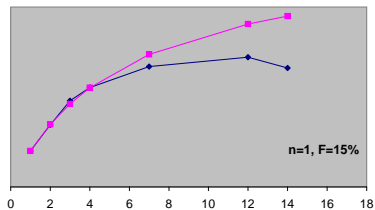
Nonlinear Estimation Experiment

- ▶ simulate measurements from the solution specified before
- ▶ generate 200 MH draws (starting from almost true values)
- ▶ use UKF Filter
- ▶ 32 parallel sequences
- ▶ for models $N = 1, N = 2$
- ▶ set thread pool size to 1, 2, 3, 4, 7, 12, 14
- ▶ all experiments use same seed

Nonlin. Est.: Computing Times vs. # Processors



Nonlin. Est.: Speedup vs. Amdahl's Law



Conclusion

- ▶ we managed to find parallel algorithms that can provide a significant speedup on multi-core computers
- ▶ it scales well up to 4-8 processors, after that the speedup decreases more quickly than the theoretical bound
- ▶ as expected, the speedup is larger in bigger models
- ▶ perfect for dual/quad-core desktops/laptops
- ▶ no penalty in sequential use
- ▶ runs everywhere without any additional configuration
- ▶ the number of processors can be adjusted during runtime due to flexible thread pool pattern

Merci de votre attention.

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