POLICY FUNCTION ITERATION, PARALLEL PROCESSING, & AN INTRODUCTION TO MEX

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BENCHMARK MODEL

Representative household chooses $\{c_t, k_{t+1}\}_{t=0}^{\infty}$ to maximize expected lifetime utility,

$$E_t \sum_{t=0}^{\infty} \beta^t \frac{c_t^{1-\sigma}}{1-\sigma},$$

subject to

$$c_t + k_{t+1} = z_t k_t^{\alpha} + (1 - \delta) k_t$$

 $z_{t+1} = (1 - \rho) \bar{z} + \rho z_t + \varepsilon_{t+1},$

where k_0, z_0 are given.

DISCRETIZED STATE SPACE

- State variables: k_t , z_t
- Number of grid points: N_k , N_z
- Given grid boundaries: $[k_{\min} \ k_{\max}]$ and $[z_{\min} \ z_{\max}]$, set grids
- For evenly spaced grids use:

$$x_{grid} = linspace(x_{min}, x_{max}, N_x), \quad x \in \{k, z\}$$

- State space: contains $N = N_k \times N_z$ independent nodes
- To create an array for each state variable where every position is a unique permutation of the state space use:

$$[k_{gr} \; z_{gr}] = \mathtt{ndgrid}(k_{grid}, z_{grid})$$

FUNCTIONAL APPROXIMATION

 Goal: Find a function that maps the state space to the policy function

$$c_t = \underbrace{c(k_t, z_t)}_{\text{True RE Solution}} \approx \underbrace{\mathcal{P}(k_t, z_t, \eta)}_{\text{Approximating Function}}$$

where η is matrix of polynomial coefficients with M elements

- Collocation Method: M = N
- Galerkin Method: M < N

FIXED-POINT ITERATION

- 1. Given $\mathcal{P}(k_t, z_t, \eta^{q-1})$, solve for k_{t+1} using the budget constraint
- 2. Given the updated state, evaluate $\mathcal{P}(k_{t+1}, z_{t+1}, \eta^{q-1})$
- 3. Use Euler equation to back out the implied policy function
- 4. Use a routine to minimize the approximating function errors over the entire state space (implies $\hat{\eta}^q$)
- 5. Set $\eta^q = (1 \lambda)\eta^{q-1} + \lambda \hat{\eta}^q$, $\lambda \in [0, 1]$
- 6. Follow iteration procedure until convergence

TIME ITERATION

- 1. Use $\mathcal{P}(k_t, z_t, \eta^{q-1})$ as a guess for the current policy function, c_t
- 2. Solve for k_{t+1} using the budget constraint
- 3. Given the updated state, evaluate $\mathcal{P}(k_{t+1}, z_{t+1}, \eta^{q-1})$
- 4. Use a nonlinear solver so c_t satisfies the equilibrium system
- 5. Update approximating function
- 6. Follow iteration procedure until convergence

GLOBAL APPROXIMATION: MONOMIAL BASIS

- Choose the set of monomials corresponding to an n^{th} order Taylor approximation of the policy function
- Complete set of basis functions is:

$$\mathcal{B}_n^s = \left\{ \left(x_1^{k_1} \cdots x_s^{k_s} \right) \middle| \sum_{i=1}^s k_i = j, k_i \ge 0, j = 0, 1, \dots, n \right\},$$

where *s* is the number of state variables

- The monomials are evaluated at each of the N nodes in the discretized state-space and stored column-wise in X
- Obtain updated coefficients using OLS:

$$\hat{\eta}^q = (X'X)^{-1}X'\Lambda_t,$$

where Λ_t is a vector of policy function values.

MONOMIAL BASIS EXAMPLE

If n = 2, 2 state variables (k and z) imply

Complete Basis:

$$\mathcal{P}(k,z,\eta) = \eta_0 + \eta_k k + \eta_z z + \eta_{kk} k^2 + \eta_{kz} kz + \eta_{zz} z^2$$

Regressor matrix (subscripts denote grid indices)

$$X = \begin{bmatrix} 1 & k_1 & z_1 & k_1^2 & k_1 z_1 & z_1^2 \\ 1 & k_2 & z_2 & k_2^2 & k_2 z_2 & z_2^2 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \vdots \\ 1 & k_N & z_N & k_N^2 & k_N z_N & z_N^2 \end{bmatrix}$$

Obtain updated coefficients using OLS:

$$\hat{\eta}_c^q = (X'X)^{-1}X'\mathbf{c},$$

LOCAL APPROXIMATION

- Linear Interpolation: 2 state variables (k, z)
- Goal: Find the policy value c(k', z')
- · We have policy function values on nearest nodes,

$$[c(k_i, z_j), c(k_i, z_{j+1}), c(k_{i+1}, z_j), c(k_{i+1}, z_{j+1})],$$

once we determine the grid indices, i, j

• Locate grid point to left of k' (evenly spaced nodes)

$$\begin{aligned} \text{step} &= k_2 - k_1, & \text{dist} &= k' - k_1 \\ \text{loc} &= \min(N_k - 1, \max(1, \text{floor}(\text{dist/step}) + 1)) \\ &\xrightarrow{\text{dist}} & \text{loc} \end{aligned}$$

LOCAL APPROXIMATION

Interpolate in the k direction to obtain

$$\begin{split} c(k',z_j) &= c(k_i,z_j) + (k'-k_i) \frac{c(k_{i+1},z_j) - c(k_i,z_j)}{k_{i+1} - k_i} \\ &= \underbrace{\frac{k_{i+1} - k'}{k_{i+1} - k_i}}_{\omega_{k_i}} c(k_i,z_j) + \underbrace{\frac{k' - k_i}{k_{i+1} - k_i}}_{\omega_{k_{i+1}}} c(k_{i+1},z_j) \end{split}$$

Then interpolate in the z direction to obtain

$$c(k',z') = \underbrace{\frac{z_{j+1} - z'}{z_{j+1} - z_j}}_{\omega_{z_j}} c(k',z_j) + \underbrace{\frac{z' - z_j}{z_{j+1} - z_j}}_{\omega_{z_{j+1}}} c(k',z_{j+1}).$$

Combine these two equations

$$c(k', z') = \sum_{a=0}^{1} \sum_{b=0}^{1} \omega_{k_{i+a}} \omega_{z_{j+b}} c(k_{i+a}, z_{j+b}),$$

FORTRAN/MEX ADVANTAGE

Use a nested loop to calculate the nested sum

- Number of nested loops corresponds to the number of state variables
- Since these are unavoidable, compiled code is faster

INTEGRATION: TRAPEZOID RULE

$$E[\Phi(\cdot, z)] \approx \frac{\Pr(\varepsilon_{1})\Phi(\cdot, z(\varepsilon_{1})) + \Pr(\varepsilon_{2})\Phi(\cdot, z(\varepsilon_{2}))}{2} \Delta \varepsilon$$

$$+ \frac{\Pr(\varepsilon_{2})\Phi(\cdot, z(\varepsilon_{2})) + \Pr(\varepsilon_{3})\Phi(\cdot, z(\varepsilon_{3}))}{2} \Delta \varepsilon + \cdots$$

$$+ \frac{\Pr(\varepsilon_{m-1})\Phi(\cdot, z(\varepsilon_{m-1})) + \Pr(\varepsilon_{m})\Phi(\cdot, z(\varepsilon_{m}))}{2} \Delta \varepsilon$$

$$= \frac{\Delta \varepsilon}{2} \left[2 \sum_{i=1}^{m} \Pr(\varepsilon_{i})\Phi(\cdot, z(\varepsilon_{i})) - \Pr(\varepsilon_{1})\Phi(\cdot, z(\varepsilon_{1})) - \Pr(\varepsilon_{m})\Phi(\cdot, z(\varepsilon_{m})) \right]$$

$$\Pr(\varepsilon_{4})\Phi(\cdot, z(\varepsilon_{4}))$$

$$\Pr(\varepsilon_{4})\Phi(\cdot, z(\varepsilon_{4}))$$

HTER AND THROCKMORTON. POLICY FUNCTION ITERATION

INTEGRATION: GAUSS-HERMITE

• Given shock, $\varepsilon \sim N(\mu, \sigma^2)$,

$$E[\Phi(\cdot, z(\varepsilon))] = (2\pi\sigma^2)^{-1/2} \int_{-\infty}^{\infty} \Phi(\cdot, z(\epsilon)) e^{-(\varepsilon - \mu)^2/(2\sigma^2)} d\epsilon$$

• Apply change of variables, $u = (\epsilon - \mu)/(\sqrt{2}\sigma)$,

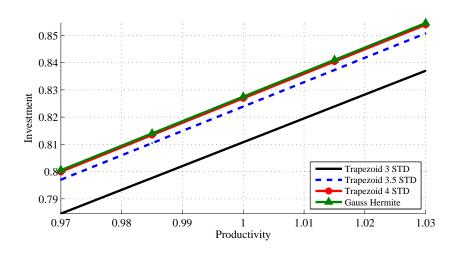
$$E[\Phi(\cdot, z(u))] = \pi^{-1/2} \int_{-\infty}^{\infty} \Phi(\cdot, z(\sqrt{2}\sigma u + \mu)) e^{-u^2} du$$
$$\approx \pi^{-1/2} \sum_{i=1}^{n} \omega_i \Phi(\cdot, z(\sqrt{2}\sigma u_i + \mu)),$$

• ω_i are Gauss-Hermite weights

$$\omega_i = 2^{n+1} n! \sqrt{\pi} [H_{n+1}(u_i)]^{-2}$$

 H_{n+1} is the physicists' Hermite polynomial of order n+1.

INTEGRATION COMPARISON



OBTAINING A GUESS

Find log-linear solution on the state space as a guess for c(k, z)

Log-linear solution from gensys takes the form,

$$\hat{Y}_{t+1} = T\hat{Y}_t + M\varepsilon_{t+1},$$

where $\hat{Y}_t = [\hat{k}_t, \hat{z}_t, \hat{c}_{t-1}]'$ and $\varepsilon \sim N(0, \sigma^2)$.

- Convert state space in levels to percent deviations from steady state
- Compute solution for i = 1, ..., N

$$\hat{c}_i = T(c_{loc}, [k_{loc} \ z_{loc}])[\hat{k}_{gr}(i) \ \hat{z}_{gr}(i)]'$$

Convert ĉ to levels

TIME ITERATION EXAMPLE

Solve for an updated c(k, z) that satisfies equilibrium system

- 1. Use log-linear solution as a guess on each node
 - If global approximation, then calculate $\hat{\eta}^0 = (X'X)^{-1}X'\mathbf{c}$
- 2. Solve for k_{t+1}, z_{t+1}
- 3. Find c_{t+1} given the updated state
 - If global approximation, update the basis. Then $c_{t+1} = X_{t+1} \eta^{q-1}$
 - · If local approximation, then use linear interpolation
- 4. Evaluate the expectation (Trapezoid rule or GH), where

$$\Phi(\varepsilon_i) = \beta(c_{t+1}(\varepsilon_i))^{-\sigma} (\alpha z_{t+1}(\varepsilon_i) k_{t+1}^{\alpha-1} + 1 - \delta)$$

TIME ITERATION EXAMPLE

- 5. Use nonlinear solver to find a c(k,z) that satisfies the consumption Euler equation, $c_t^{-\sigma}=E[\Phi]$.
- 6. Update policy function
 - Global: $\hat{\eta}^q = (X'X)^{-1}X'\mathbf{c}$.
 - · Local: use outputted policy function
- 7. Calculate distance between updates
 - Global: dist = $\eta^q \eta^{q-1}$
 - Local: dist = $c^q(k, z) c^{q-1}(k, z)$
- 8. If |dist| < tol, then stop. If not, repeat algorithm using the new policy function as the starting guess.

Advantages: Satisfies the equilibrium system on each node and nodes can be run in parallel.

Disadvantage: Non-linear solver must execute on each node.

FIXED-POINT ITERATION EXAMPLE

Solve for an updated c(k, z) implied by the equilibrium system

- 1. Use log-linear solution as a guess on each node
 - If global approximation, then calculate $\hat{\eta}^0 = (X'X)^{-1}X'\mathbf{c}$. Calculate $c(k,z) = X\hat{\eta}^0$.
- 2. Calculate updated variables and the expectation (steps 2-4 in the time iteration algorithm)
- 3. Solve for the implied policy, $c_t = (E[\Phi])^{-1/\sigma}$, on each node
- 4. Execute steps 6-8 in the time iteration algorithm

Advantage: Does not require a loop, since all of the nodes are evaluated simultaneously.

<u>Disadvantage</u>: Since resulting c(k,z) does not satisfy equilibrium system, algorithm is less stable.

EXTENSION 1: ELASTIC LABOR SUPPLY

Household now chooses $\{c_t, n_t, k_{t+1}\}_{t=0}^{\infty}$ to maximize:

$$E_t \sum_{t=0}^{\infty} \beta^t \left\{ \frac{c_t^{1-\sigma}}{1-\sigma} - \chi \frac{n_t^{1+\eta}}{1+\eta} \right\},\,$$

subject to

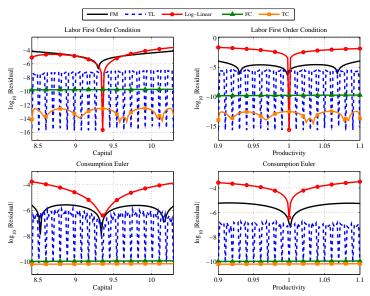
$$c_t + k_{t+1} = z_t k_t^{\alpha} n_t^{1-\alpha} + (1 - \delta) k_t$$
$$z_{t+1} = (1 - \rho) \bar{z} + \rho z_t + \varepsilon_{t+1}$$

Optimality conditions now includes:

$$(1 - \alpha)y_t/n_t = \chi n_t^{\eta} c_t^{\sigma}$$

New optimality condition is a static relation and does not add a state variable, but may require a new policy function: $n(k_t, z_t)$

EULER EQUATION ERRORS



RICHTER AND THROCKMORTON: POLICY FUNCTION ITERATION

EXTENSION 2: PRODUCTIVITY SWITCHING

Productivity evolves according to:

$$z_{t+1} = (1 - \rho)\bar{z}(s_{t+1}) + \rho z_t + \varepsilon_t,$$

where $s_t \in \{1, 2\}$, z(1) < z(2), and

$$\begin{bmatrix} \Pr[s_{t+1} = 1 | s_t = 1] & \Pr[s_{t+1} = 2 | s_t = 1] \\ \Pr[s_{t+1} = 2 | s_t = 1] & \Pr[s_{t+1} = 2 | s_t = 2] \end{bmatrix} = \begin{bmatrix} p_{11} & p_{12} \\ p_{21} & p_{22} \end{bmatrix},$$

where $0 \leq p_{ij} \leq 1$ and $\sum_{j=1}^{2} p_{ij} = 1$ for all $i \in \{1,2\}$

Key Changes:

- Policy functions are dependent on a discrete state variable
- Policy functions account for expectational effects of alternative regimes

SOLVING THE MODEL

- Calculate $z_{t+1}(s_{t+1})$ for each realization of the productivity state
- Find $c_{t+1}(s_{t+1})$ and $n_{t+1}(s_{t+1})$ based on each realization of the productivity state
- Numerical Integration:
 - First integrate across the continuous random variable, ε , to obtain expectations, conditional on the realizations of the discrete stochastic variable at time t+1. This yields:

$$E[\Phi|s_{t+1}=1], \quad E[\Phi|s_{t+1}=2]$$

 Then weight each of these values by their corresponding likelihood. Conditional expectations are:

$$E_t [\Phi | s_t = i] = p_{i1} E[\Phi | s_{t+1} = 1] + p_{i2} E[\Phi | s_{t+1} = 2]$$

Introduction to MEX

- Advantages of MATLAB:
 - Many built-in functions with good documentation
 - Parallel computing toolbox (PCT) makes parallel processing easy to implement
 - · Very easy to debug code
- Main drawback: Interpretive language is very slow at evaluating loops
- MEX (MATLAB Executable) functions: Allow programmers to write sections of the code using a compiled language and import it into MATLAB
- Challenge: Users must write a "Gateway" function that allows MATLAB to communicate with compiled code

FORTRAN 90 MEX REQUIREMENTS

- Intel Visual Fortran Composer XE 2011
 - · Basic compiler: very few intrinsic functions
 - IMSL Library: Provides hundreds of additional functions
- Microsoft Visual Studio 2008
- MATLAB default: Fixed-format (f77) Fortran code.
- Our code: Free-format (f90), which is similar to MATLAB.
- To change the default settings, modify the batch files (...\bin\win64\mexopts) by deleting the '/fixed' flag
- Use mex -setup to select IVF as the compiler in MATLAB

WRITING A GATEWAY SUBROUTINE

- Every Fortran MEX-file gateway routine begins like this: 1 #include "fintrf.h"
- 2 subroutine mexFunction(nlhs,plhs,nrhs,prhs)
 - ! Declarations
- 3 implicit none
 - ! mexFunction arguments
- 4 mwPointer plhs(*), prhs(*)
- 5 integer * 4 nlhs, nrhs
- Line 1 defines pointer types in the MATLAB interface
- Function arguments:
 - prhs: Pointer to an array which holds the input data
 - plhs: Pointer to an array which will hold the output data
 - nrhs: number of right-hand (input) arguments
 - nlhs: number of left-hand (output) arguments
- Line 3 avoids Fortran's implicit type definitions

KEY MATLAB INTERFACE FUNCTIONS

- mxGetPr: Accesses the real data in an mxArray
- mxGetScalar: Grabs the value of the first real element of the mxArray (often one element)
- mxGetM/mxGetN: Determines the number of rows/columns in a specified mxArray
- mxGetDimensions: Determines how many elements are in each dimension of the mxArray
- mxClassIDFromClassName: Obtains an identifier for any MATLAB class (e.g. Double)
- mxCreateNumericArray: Creates an N-dimensional mxArray in which all data elements have the numeric data type specified by ClassID.
- mxCreateDoubleMatrix: Creates an m-by-n mxArray

INPUTTING A POLICY FUNCTION

 Declare variable types and sizes (lines 1-3). If the dimension lengths are variable, use allocatable memory:

```
1 mwpointer c_pr
2 mwSize nk,nz
3 real*8, allocatable, dimension(:,:) :: c
! Load Inputs
4 nk = mxGetN(prhs(1))
5 nz = mxGetN(prhs(2))
6 allocate(c(nk,nz))
7 c_pr = mxGetPr(prhs(3))
8 call mxCopyPtrToReal8(c_pr,c,nk*nz)
Load the dimensions of ref from inputs (lines 4 and 5) a
```

- Load the dimensions of pf from inputs (lines 4 and 5) and allocate the memory (line 6)
- Grab the address of the pf (input 3), store in c_pr (line 7), and copy to Fortran variable c (line 8)

CREATE OUTPUT MATRIX

```
    Load output size: stochastic realizations (lines 1-2)

1 e pr = mxGetDimensions(prhs(4))
2 call mxCopyPtrToInteger4(e pr,e,1)
  !Create array for return argument
3 cid = mxClassIDFromClassName('double')
4 allocate(o(e))
5 plhs(1) = mxCreateNumericArray(1,e,cid,0)
6 o_pr = mxGetPr(plhs(1))
  ! Call subroutine and load Fortran array
7 call interpfnc(required inputs)
8 call mxCopyReal8ToPtr(o,o pr,e)
```

- Create 1 × e output vector of type double (lines 3-5) and assign address (line 6)
- Call interpolation subroutine (line 7) and copy the data to the output address (line 8)

INTRODUCTION TO PARALLEL PROCESSING

- Any calculations that are not dependent on the results of other calculations can be performed in parallel (e.g. solving for policy values at each node in the state space)
- With the PCT, a maximum of 12 processors may be used locally with any version of MATLAB 2011a or later (8 processors with 2009a or later, 4 processors with 2007a or later). Additional processors require the MATLAB Distributed Computing Server (MDCS).
- All available processors are initialized with the matlabpool command (use matlabpool open X for a subset of the available processors) and terminated with matlabpool close

INTRODUCTION TO PARALLEL PROCESSING

The PCT requires the following alterations to the code:

- Replace for loops with parfor loops where applicable.
 This tells MATLAB to distribute each step in the loop across the available processors
- If a nested for loop is used to update policy functions across dimensions, reduce it to a single loop by changing to a single index (as opposed to specifying coordinates)
- Remove all global variables and instead use structures/parameter lists and variable arrays as direct inputs into functions called within the parfor loop

ADVANTAGE OF MEX AND PARALLELIZATION

RBC Solution Times–MEX, Parallelization Comparison* (in seconds)

	MATLAB		MEX	
Processors	Structures	No Structures	Structures	No Structures
1	317 (1.0)	254 (1.2)	191 (1.7)	132 (2.4)
2	164 (1.9)	132 (2.4)	101 (3.1)	71 (4.5)
6	65 (4.9)	53 (6.0)	41 (7.8)	32 (10.0)

^{*} Based on a state space comprised of 1,681 nodes (41 points on k_t and z_t). We specify 10 realizations of ε_{t+1} . The routines were computed with an Intel W3670 6-core processor (3.20GHz) operating 64-bit Windows 7.

CONVENTIONAL NEW KEYNESIAN MODEL

The representative household chooses $\{c_{t+i}, n_{t+i}, M_{t+i}, k_{t+i}\}_{i=0}^{\infty}$ to maximize expected lifetime utility,

$$E_t \sum_{i=0}^{\infty} \beta^i \left\{ \frac{c_{t+i}^{1-\sigma}}{1-\sigma} - \chi \frac{n_{t+i}^{1+\eta}}{1+\eta} + \nu \frac{\left(M_{t+i}/P_{t+i}\right)^{1-\kappa}}{1-\kappa} \right\}, \qquad \chi, \nu > 0$$

subject to

$$c_t + m_t + i_t + b_t = (1 - \tau_t)(w_t n_t + r_t^k k_{t-1}) + (m_{t-1} + r_{t-1} b_{t-1})/\pi_t + d_t$$
$$k_t = (1 - \delta)k_{t-1} + i_t$$

Optimality implies

$$\chi n_t^{\eta} c_t^{\sigma} = (1 - \tau_t) w_t \tag{FOC n}$$

$$\nu m_t^{-\kappa} = (1 - 1/r_t) c_t^{-\sigma} \tag{FOC m}$$

$$1 = \beta r_t E_t \{ (c_t/c_{t+1})^{\sigma}/\pi_{t+1} \}$$
 (FOC b)

$$1 = \beta E_t \left\{ (c_t/c_{t+1})^{\sigma} \left(r_{t+1}^k + 1 - \delta \right) \right\}$$
 (FOC k)

CONVENTIONAL NEW KEYNESIAN MODEL

Firm optimality conditions:

$$\begin{aligned} y_t &= k_{t-1}^{\alpha} n_t^{1-\alpha} \\ r_t^k &= \alpha \Psi_t y_t / k_{t-1} \\ w_t &= (1-\alpha) \Psi_t y_t / n_t \\ \varphi\left(\frac{\pi_t}{\bar{\pi}} - 1\right) \frac{\pi_t}{\bar{\pi}} &= (1-\theta) + \theta \Psi_t + \varphi E_t \left[q_{t,t+1} \left(\frac{\pi_{t+1}}{\bar{\pi}} - 1\right) \frac{\pi_{t+1}}{\bar{\pi}} \frac{Y_{t+1}}{Y_t} \right] \end{aligned}$$

Government Budget Constraint:

$$m_t + b_t + \tau_t(w_t n_t + r_t^k k_{t-1}) = \bar{g} + \underbrace{(m_{t-1} + r_{t-1} b_{t-1})}_{\equiv a_t} / \pi_t$$

Monetary and Fiscal Policy Rules:

$$r_t = \bar{r}(\pi_t/\pi^*)^{\phi(s_t)} \exp(\varepsilon_{r,t}), \qquad \tau_t = \bar{\tau}(b_{t-1}/b^*)^{\gamma(s_t)} \exp(\varepsilon_{\tau,t})$$

Aggregate Resource Constraint:

$$c_t + i_t + \bar{q} = [1 - \varphi(\pi_t/\bar{\pi} - 1)^2/2]y_t$$

SOLVING THE MODEL: GUESS

- Minimum State Nonlinear Model: a_t , b_{t-1} , k_{t-1} , $\varepsilon_{r,t}$, $\varepsilon_{\tau,t}$, s_t
- Minimum State Linear Model: m_{t-1} , r_{t-1} , b_{t-1} , and k_{t-1} . Must average across one of the components of $a_t = m_{t-1} + r_{t-1}b_{t-1}$ to make these sets consistent and replicate across the shocks
- Use a convex combination of the linear solution for each policy mix with weights equal to the transition probabilities to form a guess for the nonlinear policy functions
 - Weighting the linear solutions approximates the expectational effect of switching between regimes

SOLVING THE MODEL: POLICIES

• Policy Functions: c_t , n_t , π_t , Ψ_t . Note: Ψ_t can be omitted if the firm's optimality conditions are rewritten as

$$\frac{k_{t-1}}{n_t} = \frac{\alpha}{1 - \alpha} \frac{w_t}{r_t^k}$$

$$\Psi_t = w_t^{1-\alpha} (r_t^k)^{\alpha} (1 - \alpha)^{-(1-\alpha)} \alpha^{-\alpha}$$

- Given the state and policies, solve for all time t variables
- Interpolate values for the updated policies across both shocks and solve for variables inside expectations
 - With two continuous stochastic components, interpolated policies have $N_{\varepsilon_{\tau}} \times N_{\varepsilon_{\tau}}$ realizations (a matrix of values). Requires a nested loop in the interpolation function.

SOLVING THE MODEL: INTEGRATION

- Must numerically integrate across both random variables:
 - Replicate $\Pr(\varepsilon_{\tau,i})$ (the row weights) across the $\varepsilon_{r,j}$ -dimension (the columns) to create an $N_{\varepsilon_{\tau}} \times N_{\varepsilon_{r}}$ weighting matrix
 - Use the weighting matrix to weight each element of $[\Phi_{i,j}]$
 - Use either the Trapezoid rule or GH to integrate across the ε_{τ} -dimension (the rows)—collapses the matrix of realizations to a row vector
 - Weight each of the outcomes in the row vector by its respective $\Pr(\varepsilon_{r,j})$
 - Use either the Trapezoid rule or GH to integrate across the ε_r -dimension—reduces to a single value (state dependent)
- Use transition probabilities to integrate across the discrete random variable (policy regime state)

ADVANTAGE OF MEX AND PARALLELIZATION

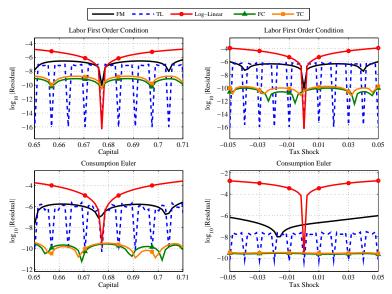
 \star Runs are based on 1 continuous stochastic variable (ε_{τ}) and fixed AM/PF policy

NK Model Solution Times* (in seconds)

Processors	No MEX	Partial MEX	All MEX
1	288.2 (1.0)	131.5 (2.2)	32.0 (9.0)
2	148.7 (1.9)	68.6 (4.2)	16.2 (16.2)
6	54.1 (5.3)	27.0 (10.7)	8.2 (35.1)

^{*} Based on a state space comprised of 2,401 nodes (7 points on each continuous state variable). We specify 10 realizations of ε_{τ} . The routines were computed with an Intel W3670 6-core processor (3.20GHz) operating 64-bit Windows 7. The value in parenthesis represents that speed factor increase from the slowest setup.

EULER EQUATION ERRORS (FIXED AM/PF)



SPEED COMPARISON

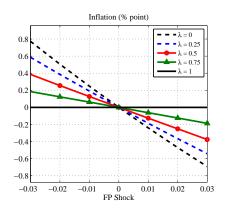
Solution Times-Solution Methods Comparison* (in seconds)

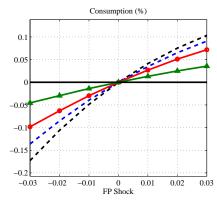
	RBC		NK	
Method	MATLAB	MEX	MATLAB	MEX
FM	3	N/A	3	N/A
FC	139	6	948	19
TL	62	42	69	39
TC	106	44	258	40

^{*} RBC solution is based on a state space comprised of 1,681 nodes (41 points on k_{t-1} , 41 points on z_t). We specify 10 realizations of ε_{t+1} . NK solution is based on a state space comprised of 2,401 nodes (7 points on $a_{t-1}, b_{t-1}, k_{t-1}$, and $\varepsilon_{\tau,t}$). We specify 10 realizations of $\varepsilon_{\tau,t+1}$. The routines were computed with an Intel W3670 6-core processor (3.20GHz) operating 64-bit Windows 7. All locally available processors were used.

POLICY FUNCTIONS (LUMP-SUM TAXES)

- Average duration of time spent in the AM/PF regime in the ergodic distribution: $\lambda=(1-p_{22})/(2-p_{11}-p_{22})$
- When $\lambda = 1$, there is no chance of moving to the PM/AF

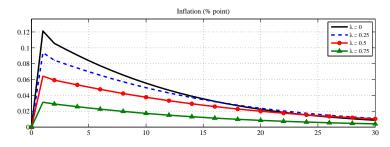


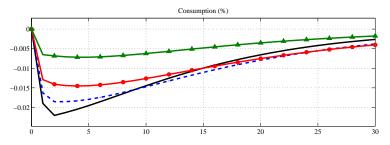


IMPULSE RESPONSES

- Nonlinear impulse responses differ from linear responses
- Linear impulse responses report how a shock makes each simulated variable differ from its deterministic (calculated) steady state
- Nonlinear impulse responses report how a shock makes each simulated path differ from the *stochastic* steady state, defined as a value for the state vector such that $|\Theta_t \Theta_{t-1}| < \epsilon$, where Θ is a vector of policy functions
- To simulate the model, burn off the number of periods needed to ensure that the model reaches its stochastic steady state
- Although the transversality condition is not imposed, simulated paths in models that explicitly violate this condition will diverge even if the algorithm converges

IMPULSE RESPONSE: FP SHOCK





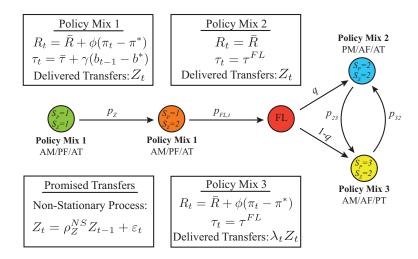
RICHTER AND THROCKMORTON: POLICY FUNCTION ITERATION

EXPLOSIVE TRANSFERS AND A FISCAL LIMIT

- State: a_t , b_{t-1} , k_{t-1} , z_t , $s_{z,t}$, $s_{p,t}$ (no shocks on MP or FP)
- z: lump-sum transfers to household, distributed AR(1)
 - $s_z = 1$: transfers process is stationary.
 - $s_z = 2$: transfers process is non-stationary.
- s_p: monetary/fiscal policy mix
 - $s_p = 1$: active MP, passive FP, and promised transfers (Before Fiscal Limit)
 - $s_p = 2$: passive MP, active FP, and promised transfers (After Fiscal Limit)
 - $s_p = 3$: active MP, active FP, and delivered transfers (After Fiscal Limit)
- The probability of hitting the fiscal limit—point where higher taxes are no longer feasible—is endogenous

$$p_{FL,t} = 1 - \frac{\exp(\eta_0 - \eta_1(\tau_{t-1} - \bar{\tau}))}{1 + \exp(\eta_0 - \eta_1(\tau_{t-1} - \bar{\tau}))}$$

Possible Policy Evolution



OBTAINING A GUESS

To obtain initial conjectures for the policy functions:

- First solve the *linear* model under both the initial (AM/PF/AT) and debt revaluation (PM/AT/AT) regimes
- Use these linear solutions as initial guesses for the corresponding fixed-regime non-linear models, assuming there is no probability of hitting the fiscal limit
- Use weighted averages of these non-linear solutions to obtain initial conjectures for each regime combination (3 policy states and 2 transfers states form 6 combinations)

SOLVING THE MODEL

- Given the state and numerical policies, calculate time t variables (state dependent)
- Calculate updated (time t+1) values for each of the policy variables using piecewise linear interpolation/extrapolation. Note that s_z and s_p form 6 possible outcomes.
- Solve for the variables that are inside expectations (obtain all other t+1 values from the interpolated policies)
- Integrate across the continuous transfers shock. Each expectation will be conditional on the 6 possible outcomes
- Calculate p_{FL} and integrate across the policy regimes (s_p)
- Integrate across the transfers regimes (s_z)
- Using Chris Sims' root finder, csolve, solve for the zeros
 of the equations with embedded expectations, subject to
 each of the remaining equilibrium conditions.

STABILITY AND GRID ADJUSTMENTS

- Achieve stability by starting with a narrow grid near the initial steady state
- · Simulate the model and identify areas of extrapolation
- Since non-stationary transfers cause drift, must make small adjustments to the grid using previous nonlinear solution as guess
- Continue adjusting the grids until they encompass all simulated paths of state variables

MONTE CARLO SIMULATIONS

- Pick a large number of times to simulate the model
- Begin all simulations from stochastic steady state
- Take draws from the distributions of the random variables using random number generators
 - $\varepsilon_{z,t} \sim i.i.d.N(0,\sigma_z^2)$
 - For Markov processes, draw from U(0,1) and create sequences for each discrete state variable (transfers and policy regimes) using the respective transition matrix
- Use the quantile function to calculate paths that characterize the tails of the distribution for variables of interest across time, e.g., 10th and 90th percentile of all realizations