## Numerical Methods for New Keynesian ZLB Models

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#### What is covered

- Equilibrium with a simple Taylor rule
- Tauchen's (1986) method for approximating AR(1) process
- Equilibrium under the optimal discretionary policy (Adam and Billi, 2007)

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#### Two-state shock process

• Exogenous shocks take only  $N_s=2$  values,  $s_t \in \{s_H, s_L\}$ . The stochastic process follows a Markov chain with the transition matrix:

$$\left[\begin{array}{cc} 1-p_H & p_H \\ 1-p_L & p_L \end{array}\right].$$

ullet  $p_H$  is the frequency of crisis and  $p_L$  is the duration of crisis.

#### Equilibrium with Taylor rule

• Equilibrium conditions are

$$y_{t} = E_{t}y_{t+1} - (i_{t} - E_{t}\pi_{t+1} - s_{t}),$$
  

$$\pi_{t} = \kappa y_{t} + \beta E_{t}\pi_{t+1},$$
  

$$i_{t}^{*} = r^{*} + \phi_{\pi} E_{t}\pi_{t+1},$$

and the zero lower bound

$$i_t = \{0, i_t^*\}.$$

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## Analytical solution

• The solution has a form of (we omit time subscripts for the policy function)

$$y = y(s), \quad \pi = \pi(s), \quad i = i(s).$$

We know that the functions have only two values, i.e.,

$$y = \begin{cases} y_H, & \pi = \begin{cases} \pi_H, \\ \pi_L, \end{cases} \quad i = \begin{cases} i_H, \\ i_L. \end{cases}$$

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# Analytical solution, cont'd

• We assume that  $i_H > 0$  and  $i_L = 0$ . Then we have

$$\begin{split} y_H &= (1-p_H)y_H + p_H y_L - (i_H - [(1-p_H)\pi_H + p_H \pi_L] - s_H) \,, \\ \pi_H &= \kappa y_H + \beta \left[ (1-p_H)\pi_H + p_H \pi_L \right] \,, \\ i_H &= r^* + \phi_\pi \left[ (1-p_H)\pi_H + p_H \pi_L \right] \,, \\ y_L &= (1-p_L)y_H + p_L y_L - \left( 0 - [(1-p_L)\pi_H + p_L \pi_L] - s_L \right) \,, \\ \pi_L &= \kappa y_L + \beta \left[ (1-p_L)\pi_H + p_L \pi_L \right] \,, \\ i_L &= 0 \,. \end{split}$$

• There are 6 equations and 6 unknowns, so we can solve for the unknowns.

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# Numerical solution: Policy function iteration

• Consider the general case of  $N_s \geq 2$ . A guess of the policy functions

$$y = y^{(0)}(s), \quad \pi = \pi^{(0)}(s), \quad i = i^{(0)}(s).$$

• We know the values of the functions only at each grid point, e.g.,

$$\begin{split} y^{(0)}(s) &= [y_1, y_2, ..., y_{N_s}]', \\ \pi^{(0)}(s) &= [\pi_1, \pi_2, ..., \pi_{N_s}]', \\ i^{(0)}(s) &= [i_1, i_2, ..., i_{N_s}]'. \end{split}$$

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• At each grid point  $k = 1, ..., N_s$ , we solve

$$y_k = y^e - (i_k - \pi^e - s_k),$$
  
 $\pi_k = \kappa y_k + \beta \pi^e,$   
 $i_k = \max\{0, r^* + \phi_\pi \pi^e\},$ 

for  $(y_k, \pi_k, i_k)$ , where

$$y^{e} = \sum_{l=1}^{N_{s}} p(k, l) y^{(0)}(s_{l}),$$
$$\pi^{e} = \sum_{l=1}^{N_{s}} p(k, l) \pi^{(0)}(s_{l}).$$

and p(k, l) is the (k, l) element of the transition matrix.



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• Once this is done for all the grid points, we update

$$y^{(1)}(s) = [y_1, y_2, ..., y_N]',$$
  

$$\pi^{(1)}(s) = [\pi_1, \pi_2, ..., \pi_N]',$$
  

$$i^{(1)}(s) = [i_1, i_2, ..., i_N]'.$$

• We repeat the procedure until the policy functions converge, i.e.,  $\left\|x^{(j)}(s)-x^{(j-1)}(s)\right\|<\epsilon$  for  $x\in\{y,\pi,i\}.$ 

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#### Tauchen's method

• Tauchen (1986) developed a method for approximating AR(1) process by using Markov chain.

TBD

• The policymaker chooses  $\{\pi_t, y_t, i_t\}$  so as to maximize

$$V_0 \equiv -E_0 \sum_{t=0}^{\infty} \beta^t \left( \pi_t^2 + \lambda y_t^2 \right)$$

subject to

$$y_{t} = E_{t}y_{t+1} - (i_{t} - E_{t}\pi_{t+1}) + g_{t},$$
  

$$\pi_{t} = \kappa y_{t} + \beta E_{t}\pi_{t+1} + u_{t},$$
  

$$i_{t} \ge 0,$$

taking  $E_t y_{t+1}$  and  $E_t \pi_{t+1}$  as given.

• Exogenous shocks are given by

$$g_t = (1 - \rho_g)g + \rho_g g_{t-1} + \varepsilon_{g,t},$$
  
$$u_t = \rho_u u_{t-1} + \varepsilon_{u,t},$$

where  $\varepsilon_{g,t} \sim N(0,\sigma_g^2)$  and  $\varepsilon_{u,t} \sim N(0,\sigma_u^2)$ .

 $\bullet \ \ \mathsf{Note that} \ g = r^*.$ 

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- We know that Markov-perfect equilibrium has only natural state variables.
- Lagrangean is

$$\mathcal{L} \equiv E_0 \sum_{t} \beta^t \left( \pi_t^2 + \lambda y_t^2 \right) + 2\phi_{PC,t} \left( -\pi_t + \kappa y_t + \beta E_t \pi_{t+1} + u_t \right) + 2\phi_{EE,t} \left( -y_t - i_t + E_t y_{t+1} + E_t \pi_{t+1} + g_t \right) + 2\phi_{ZLB,t} i_t.$$

First-order necessary conditions are

$$\begin{split} \partial \pi_t : & \pi_t - \phi_{PC,t} = 0, \\ \partial y_t : & \lambda y_t + \kappa \phi_{PC,t} - \phi_{EE,t} = 0, \\ \partial i_t : & - \phi_{EE,t} + \phi_{ZLB,t} = 0. \end{split}$$

Complementary slackness condition:

$$\phi_{ZLB,t} > 0 \perp i_t > 0.$$

• When  $i_t > 0$ ,  $\phi_{ZLB,t} = 0$ . Equilibrium conitions are

$$\begin{split} i_t &= -y_t + E_t y_{t+1} + E_t \pi_{t+1} + g_t, \\ \pi_t &= \kappa y_t + \beta E_t \pi_{t+1} + u_t, \\ 0 &= \lambda y_t + \kappa \pi_t. \end{split}$$

• When  $i_t = 0$ ,  $\phi_{ZLB,t} > 0$ . Equilibrium conitions are

$$0 = -y_t + E_t y_{t+1} + E_t \pi_{t+1} + g_t,$$
  

$$\pi_t = \kappa y_t + \beta E_t \pi_{t+1} + u_t,$$
  

$$\phi_{ZLB,t} = \lambda y_t + \kappa \pi_t.$$

## Analytical solution

The solution has a form of

$$y = y(g, u), \quad \pi = \pi(g, u), \quad i = i(g, u).$$

 Now consider the case in which there are only two-state g shocks. We know that the functions have only two values, i.e.,

$$y = \begin{cases} y_H, & \pi = \begin{cases} \pi_H, \\ \pi_L, \end{cases} \quad i = \begin{cases} i_H, \\ i_L. \end{cases}$$

## Analytical solution

• We assume that  $i_H > 0$  and  $i_L = 0$ . Then we have

$$\begin{split} y_{H} &= (1-p_{H})y_{H} + p_{H}y_{L} - (i_{H} - [(1-p_{H})\pi_{H} + p_{H}\pi_{L}]) + g_{H}, \\ \pi_{H} &= \kappa y_{H} + \beta \left[ (1-p_{H})\pi_{H} + p_{H}\pi_{L} \right], \\ 0 &= \lambda y_{H} + \kappa \pi_{H}, \\ y_{L} &= (1-p_{L})y_{H} + p_{L}y_{L} - (0 - [(1-p_{L})\pi_{H} + p_{L}\pi_{L}]) + g_{L}, \\ \pi_{L} &= \kappa y_{L} + \beta \left[ (1-p_{L})\pi_{H} + p_{L}\pi_{L} \right], \\ \phi_{L} &= \lambda y_{L} + \kappa \pi_{L}. \end{split}$$

• There are 6 equations and 6 unknowns, so we can solve for the unknowns.

#### Numerical solution: Joint shock process

 Let's get back to the general case. The shock processes are appoximated by Markov chains. That is,

$$g_m \in [g_1, g_2, ..., g_{N_g}],$$
  
 $u_n \in [u_1, u_2, ..., u_{N_u}],$ 

and

$$P^g = \left[ \begin{array}{cccc} p_{11}^g & \cdots & \cdots & p_{1N_g}^g \\ p_{21}^g & \ddots & & \vdots \\ \vdots & & \ddots & \vdots \\ p_{N_g1}^g & \cdots & \cdots & p_{N_gN_g}^g \end{array} \right], \quad P^u = \left[ \begin{array}{cccc} p_{11}^u & \cdots & \cdots & p_{1N_g}^u \\ p_{21}^u & \ddots & & \vdots \\ \vdots & & \ddots & \vdots \\ p_{N_g1}^u & \cdots & \cdots & p_{N_gN_g}^u \end{array} \right].$$

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#### Joint shock process

- A kronecker product  $s = g \otimes u$  represents the joint shock process.
  - $\bullet$  For example, when  $N_g=N_u=2$

$$s_1 = (g_1, u_1),$$
  
 $s_2 = (g_1, u_2),$   
 $s_3 = (g_2, u_1),$   
 $s_4 = (g_2, u_2).$ 

- Note that each index points to a pair of shocks,  $s_k = (g_{m(k)}, u_{n(k)})$ .
- A kronecker product of the transition matrices  $P^s=P^g\otimes P^u$  is the transition matrix of the joint shock process.

## Numerical solution: Policy function iteration

• A guess of the policy functions

$$y = y^{(0)}(s), \quad \pi = \pi^{(0)}(s), \quad i = i^{(0)}(s).$$

• We know the values of the functions only at each grid point, e.g.,

$$y^{(0)}(s) = [y_1, y_2, ..., y_N]',$$
  

$$\pi^{(0)}(s) = [\pi_1, \pi_2, ..., \pi_N]',$$
  

$$i^{(0)}(s) = [i_1, i_2, ..., i_N]'.$$

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• At each grid point k = 1, ..., N, we solve

$$i_k = -y_k + y^e + \pi^e + g_{m(k)},$$
  

$$\pi_k = \kappa y_k + \beta \pi^e + u_{n(k)},$$
  

$$0 = \lambda y_k + \kappa \pi_k,$$

for  $(y_k, \pi_k, i_k)$ , where

$$y^{e} = \sum_{l=1}^{N_{s}} P^{s}(k, l) y^{(0)}(s_{l}),$$
  
$$\pi^{e} = \sum_{l=1}^{N_{s}} P^{s}(k, l) \pi^{(0)}(s_{l}).$$

• Check  $i_k \geq 0$ . If not, we solve instead

$$0 = -y_k + y^e + \pi^e + g_{m(k)},$$
  

$$\pi_k = \kappa y_k + \beta \pi^e + u_{n(k)},$$

for  $(y_k, \pi_k)$ , and set  $i_k = 0$ .

• Once this is done for all the grid points, we update

$$y^{(1)}(s) = [y_1, y_2, ..., y_N]',$$
  

$$\pi^{(1)}(s) = [\pi_1, \pi_2, ..., \pi_N]',$$
  

$$i^{(1)}(s) = [i_1, i_2, ..., i_N]'.$$

• We repeat the procedure until the policy functions converge, i.e.,  $\left\|x^{(j)}(s)-x^{(j-1)}(s)\right\|<\epsilon$  for  $x\in\{y,\pi,i\}.$ 



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