# WEEK 2: LOG-LINEARIZATION & BLANCHARD-KAHN (1980)

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# TODAY'S AGENDA

- 1. Log-linearization
- 2. Blanchard-Kahn (1980) Method based on Eric Sims Graduate Macro notes

# WHY LOG-LINEARIZE?

- Solutions to most DSGE models involve non-linear difference equations
- Log-linearization is an approximation technique that allows us to
  - 1. express the solution as a system of **linear** differential equations
  - 2. express variables in percentage terms which is easy to interpret

# **LOG-LINEARIZATION - THREE STEPS**

- 1. Log-transform all equilibrium conditions
- 2. Linearize (firt-order approx.) each equation around a point  $\rightarrow$  steady-state
- 3. Re-arrange to express everything in terms of % deviation from steady-state

# TAYLOR'S THEOREM

 An arbitrary function f(x) can be expressed as a power series about a point x\* ∈ X:

$$f(x) = f(x^*) + \frac{f'(x^*)}{1!}(x - x^*) + \frac{f''(x^*)}{2!}(x - x^*)^2 + \frac{f'''(x^*)}{3!}(x - x^*)^3 + \dots$$

 If the higher order derivatives are small, then the function can be well approximated linearly by:

$$f(x) \approx f(x^*) + \frac{f'(x^*)}{1!}(x - x^*)$$

Straightforward to extend to multi-variate functions

$$f(x,y) \approx f(x^*,y^*) + \frac{f_x'(x^*,y^*)}{1!}(x-x^*) + \frac{f_y'(x^*,y^*)}{1!}(y-y^*)$$

# **GENERIC CASE**

Consider the non-linear function:

$$f(x, y) = \frac{g(x, y)}{h(x, y)}$$

• Taking logs on both sides we get:

$$\ln f(x, y) = \ln g(x, y) - \ln h(x, y)$$

Next: compute the first order Taylor expansion for each term

# FIRST ORDER TAYLOR EXPANSION

$$\ln f(x,y) \approx \ln f(x^*,y^*) + \frac{d \ln f(x^*,y^*)}{dx^*} (x-x^*) + \frac{d \ln f(x^*,y^*)}{dy^*} (y-y^*)$$

$$\ln g(x,y) \approx \ln g(x^*,y^*) + \frac{d \ln g(x^*,y^*)}{dx^*} (x-x^*) + \frac{d \ln g(x^*,y^*)}{dy^*} (y-y^*)$$

$$\ln h(x,y) \approx \ln h(x^*,y^*) + \frac{d \ln h(x^*,y^*)}{dx^*} (x-x^*) + \frac{d \ln h(x^*,y^*)}{dy^*} (y-y^*)$$

Chain rule implies:

$$\frac{d\ln f(x)}{dx} = \frac{f'(x)}{f(x)}$$

We will use this to simplify the above expressions

## FIRST ORDER TAYLOR EXPANSION - cont.

• Combine the 3 expressions from the previous slide:

$$\ln f(x^*, y^*) + \frac{f_X'(x^*, y^*)}{f(x^*, y^*)}(x - x^*) + \frac{f_Y'(x^*, y^*)}{f(x^*, y^*)}(y - y^*) =$$

$$\ln g(x^*, y^*) + \frac{g_X'(x^*, y^*)}{g(x^*, y^*)}(x - x^*) + \frac{g_Y'(x^*, y^*)}{g(x^*, y^*)}(y - y^*) -$$

$$\left[\ln h(x^*, y^*) + \frac{h_X'(x^*, y^*)}{h(x^*, y^*)}(x - x^*) + \frac{h_Y'(x^*, y^*)}{h(x^*, y^*)}(y - y^*)\right]$$

Note that terms in pink cancel out

# FIRST ORDER TAYLOR EXPANSION - cont.

• We are left with:

$$\frac{f_{x}^{'}(x^{*},y^{*})}{f(x^{*},y^{*})}(x-x^{*}) + \frac{f_{y}^{'}(x^{*},y^{*})}{f(x^{*},y^{*})}(y-y^{*}) = \frac{g_{x}^{'}(x^{*},y^{*})}{g(x^{*},y^{*})}(x-x^{*}) + \frac{g_{y}^{'}(x^{*},y^{*})}{g(x^{*},y^{*})}(y-y^{*}) - \left[\frac{h_{x}^{'}(x^{*},y^{*})}{h(x^{*},y^{*})}(x-x^{*}) + \frac{h_{y}^{'}(x^{*},y^{*})}{h(x^{*},y^{*})}(y-y^{*})\right]$$

- Let  $\tilde{x} = \frac{x x^*}{x^*}$  be the percentage deviation of x from  $x^*$
- Multiply and divide each term by either  $x^*$  or  $y^*$

$$\frac{x^*f_x^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{x} + \frac{y^*f_y^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{y} = \frac{x^*g_x^{'}(x^*,y^*)}{g(x^*,y^*)}\tilde{x} + \frac{y^*g_y^{'}(x^*,y^*)}{g(x^*,y^*)}\tilde{y} - \left[\frac{x^*h_x^{'}(x^*,y^*)}{h(x^*,y^*)}\tilde{x} + \frac{y^*h_y^{'}(x^*,y^*)}{h(x^*,y^*)}\tilde{y}\right]$$

• The above expression is linear in both  $\tilde{x}$  and  $\tilde{y}$ 



#### LOG-LINEARIZATION - RECAP

- Take logs
- Do a first order Taylor approximation
- Multiply/divide as necessary to get variables expressed as percentage deviations from x\*
- The final product is an expression linear in  $\tilde{x}$ ,  $\tilde{y}$ :

$$\frac{x^*f_\chi^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{x} + \frac{y^*f_y^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{y} = \frac{x^*g_\chi^{'}(x^*,y^*)}{g(x^*,y^*)}\tilde{x} + \frac{y^*g_y^{'}(x^*,y^*)}{g(x^*,y^*)}\tilde{y} - \frac{x^*h_\chi^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{x} - \frac{y^*h_\chi^{'}(x^*,y^*)}{f(x^*,y^*)}\tilde{y}$$

#### LOG-LINEARIZATION - EXAMPLES

- Cobb-Douglas production function:  $y_t = a_t k_t^{\alpha} n_t^{1-\alpha}$
- Taking logs

$$\ln y_t = \ln a_t + \alpha \ln k_t + (1 - \alpha) \ln n_t$$

• First order Taylor approximation around  $y^*, a^*, k^*, n^*$ :

$$\ln y^* + \frac{1}{y^*} (y_t - y^*) = \ln a^* + \frac{1}{a^*} (a_t - a^*) + \alpha \ln k^* + \frac{\alpha}{k^*} (k_t - k^*)$$
$$+ (1 - \alpha) \ln n^* + \frac{(1 - \alpha)}{n^*} (n_t - n^*)$$

Pink terms cancel out. Re-arranging:

$$\frac{(y_t - y^*)}{y^*} = \frac{(a_t - a^*)}{a^*} + \alpha \frac{(k_t - k^*)}{k^*} + (1 - \alpha) \frac{(n_t - n^*)}{n^*}$$

• As before, we define  $\tilde{x}_t = \frac{(x_t - x^*)}{x^*}$ :

$$\tilde{y}_t = \tilde{a}_t + \alpha \tilde{k}_t + (1-\alpha) \tilde{n}_t$$



# BLANCHARD & KAHN (1980) METHOD

- After log-linearizing, we have a system of linear difference equations
- Next → how can we solve these systems?
- Four Steps
  - 1. Write model in state-space form
  - 2. Decouple the system using Jordan decomposition
  - 3. Intuition + algorithm to solve model
  - 4. Existence and uniqueness
- Dynare/Python/Julia implement this solution method
- More details (for example) here Eric Sims Notes

#### 1. WRITE MODEL IN STATE-SPACE FORM

$$\mathbb{E}_{t}[x_{t+1}] = \Psi x_{t}$$

$$x_{t} \in \mathbb{R}^{p+m}, \quad x_{t} = \left[x_{t}^{s}, x_{t}^{j}\right]$$

$$x_{t}^{s} \in \mathbb{R}^{m}, \quad x_{t}^{j} \in \mathbb{R}^{p}$$

$$x_{0}^{s} \text{ is given}$$

- $x_t$  is a vector of variables expressed in percentage deviations from ss (our  $\tilde{x}_t$ )
- We will separate this vector into two groups:
  - 1. p "jump" or forward-looking variables
  - 2. *m* state or predetermined variables
- $\Psi$  governs the evolution of the system, given a starting point
- The initial value of state variables is given (eg.  $k_0$  in the RBC model)
- **Q:** where do we start the jump variables?

## 1. WRITE MODEL IN STATE-SPACE FORM - EXAMPLE

 Log-linearized equilibrium conditions of (deterministic) neoclassical growth model:

$$-\sigma \tilde{c}_t = -\sigma \tilde{c}_{t+1} + \beta (\alpha - 1) R \tilde{k}_{t+1} \qquad \qquad \tilde{k}_{t+1} = \frac{1}{\beta} \tilde{k}_t - \frac{c}{k} \tilde{c}_t$$

Re-arranging we get:

$$\underbrace{\begin{bmatrix} \tilde{c}_{t+1} \\ \tilde{k}_{t+1} \end{bmatrix}}_{\mathbb{E}_{t}[X_{t+1}]} = \underbrace{\begin{bmatrix} 1 - \frac{c}{k} \frac{\beta(\alpha-1)R}{\sigma} & \frac{(\alpha-1)R}{\sigma} \\ -\frac{c}{k} & \frac{1}{\beta} \end{bmatrix}}_{\mathbb{\Psi}} \underbrace{\begin{bmatrix} \tilde{c}_{t} \\ \tilde{k}_{t} \end{bmatrix}}_{X_{t}}$$

- State: capital; Jump: consumption
- Note that I drop the expectation operator because the model is deterministic

# 2. DECOUPLE THE SYSTEM USING JORDAN DECOMPOSITION

• An eigenvalue is a scalar,  $\lambda_i$ , and an eigenvector is a vector  $\mathbf{v}_i$ , that jointly satisfy:  $\Psi \mathbf{v}_{\mathbb{A}} = \lambda_i \mathbf{v}_i \qquad \forall \quad i \in p+m$ 

$$\Gamma = \begin{bmatrix} v_{1,1} & v_{2,1} & \dots & v_{p+m,1} \\ \dots & \dots & \dots & \dots \\ v_{1,p+m} & v_{2,p+m} & \dots & v_{p+m,p+m} \end{bmatrix} \qquad \qquad \Lambda = diag(\lambda)$$

• We can decompose  $\Psi$  as:  $\Psi\Gamma = \Gamma\Lambda \rightarrow \Psi = \Gamma\Lambda\Gamma^{-1}$ 

# 2. DECOUPLE THE SYSTEM USING JORDAN DECOMPOSITION - II

Use Jordan decomposition to re-write model as:

$$\mathbb{E}_t[x_{t+1}] = \Gamma \Lambda \Gamma^{-1} x_t$$

• Pre-multiply each side by  $\Gamma^{-1}$ :

$$\mathbb{E}_t[\Gamma^{-1}x_{t+1}] = \Lambda \Gamma^{-1}x_t$$

• Change of variables:  $w_t = \Gamma^{-1} x_t$ . We get:

$$\mathbb{E}_{t}[w_{t+1}] = \Lambda w_{t} \to \mathbb{E}_{t} \begin{bmatrix} w_{1,t+1} \\ w_{2,t+1} \end{bmatrix} = \begin{bmatrix} \Lambda_{1} & 0 \\ 0 & \Lambda_{2} \end{bmatrix} \begin{bmatrix} w_{1,t} \\ w_{2,t} \end{bmatrix}$$

- Partition A in two:
  - 1.  $\Lambda_1$  contains Q stable eigenvalues ( $|\lambda_i| < 1$ )
  - 2.  $\Lambda_2$  contains B unstable eigenvalues ( $|\lambda_i| > 1$ )
- System is in VAR(1) form,  $\Lambda$  is diagonal  $\to w_{1,t}, w_{2,t}$  evolve independently of each other

# 2. DECOUPLE THE SYSTEM USING JORDAN DECOMPOSITION-III

Iterating forward we can write:

$$\mathbb{E}_t[w_{1,t+T}] = \Lambda_1^T w_{1,t} \qquad \qquad \mathbb{E}_t[w_{2,t+T}] = \Lambda_2^T w_{2,t}$$

- Eigenvalues in  $\Lambda_1$  are all stable  $\to$   $\Lambda_1^T \to 0$  as  $T \to \infty$
- Eigenvalues in  $\Lambda_2$  are all unstable  $\to$   $\Lambda_2^T \to \infty$  as  $T \to \infty$
- This implies  $\mathbb{E}_t[w_{2,t=T}] \to \infty$  which is inconsistent with transversality condition /bounded equilibrium unless  $w_{2,t} = 0$
- Goal: find policy function/ updating rule for jump variables by setting  $w_{2,t} = 0$

## 3. INTUITION + ALGORITHM TO SOLVE MODEL

• We can express w<sub>t</sub> as:

$$\underbrace{w_{1,t}}_{Q\times 1} = \underbrace{G_{11}}_{Q\times p} \underbrace{x_t^j}_{p\times 1} + \underbrace{G_{12}}_{Q\times m} \underbrace{x_t^s}_{m\times 1} \quad \text{where} \quad \Gamma^{-1} = \begin{bmatrix} G_{11} & G_{12} \\ G_{21} & G_{22} \end{bmatrix}$$

$$\underbrace{w_{2,t}}_{B\times 1} = \underbrace{G_{21}}_{B\times p} \underbrace{x_t^j}_{p\times 1} + \underbrace{G_{22}}_{B\times m} \underbrace{x_t^s}_{m\times 1}$$

• Set  $w_{2,t} = 0$ :

$$\underbrace{\mathbf{0}}_{\text{21}} = G_{21} x_t^j + G_{22} x_t^s \longrightarrow x_t^j = -G_{21}^{-1} G_{22} x_t^s$$

- When will  $G_{21}$   $(B \times p)$  have an inverse? Necessary condition: p = B.
- In other words, we need as many jump variables as unstable eigenvalues

## 3. INTUITION + ALGORITHM TO SOLVE MODEL

• When p = B, we get:

$$x_t^j = \underbrace{-G_{21^{-1}G_{22}}}_{P} x_t^s$$

- Recall that the starting point of the state variables,  $x_0^{S}$ , is given
- We can use the policy function to find the period-zero value of the jump variables that yields  $w_{2,0} = 0$
- Once we have  $x_0^s$  and  $x_0^j$ , we can use the expression in previous slide to get  $w_{1,0}$
- Then, we can use  $\mathbb{E}_t[w_{1,t+T}] = \Lambda_1^T w_{1,t}$  and  $\mathbb{E}_t[w_{2,t+T}] = \Lambda_2^T w_{2,t}$  to compute the evolution of the system across time

# 4. EXISTENCE AND UNIQUENESS OF A SOLUTION

- 1. If p = B, there exists a unique bounded equilibrium
- 2. If p < B (more explosive  $\lambda$ 's than jump variables), there is no solution
- 3. If p > B, there is an infinite number of solutions.