Financial Theory IV: Solving Structural Models

Dan Greenwald

Spring 2024

Dynamic Programming

- Let x_t be endogenous states, z_t be exogenous states, and y_t be controls.
- Basic problem:

$$V(x_{t}, z_{t}) = \max_{y_{t} \in \Gamma(x_{t}, z_{t})} F(x_{t}, y_{t}, z_{t}) + \beta E_{t} [V(x_{t+1}, z_{t+1})]$$
$$x_{t+1} = g(x_{t}, y_{t}, z_{t})$$
$$z_{t+1} = h(z_{t}, \varepsilon_{t+1})$$

Example: consumption-savings problem.

$$V(a_{t}, w_{t}) = \max_{x_{t} \ge 0} u (a_{t} + w_{t}\bar{L} - s_{t}) + \beta E_{t} [V(a_{t+1}, w_{t+1})]$$

$$a_{t+1} = Rs_{t}$$

$$\log w_{t+1} = (1 - \rho) \log \bar{w} + \rho \log w_{t} + \varepsilon_{t+1}, \qquad \varepsilon_{t+1} \sim N(0, \sigma^{2})$$

Dynamic Programming: Discrete Models

- ▶ Very simple and robust approach: assume $x_t \in \mathcal{X} = (\bar{x}_1, ..., \bar{x}_N)$, $z_t \in \mathcal{Z} = (\bar{z}_1, ..., \bar{z}_K)$.
 - Easy to estimate time series using Hamilton filter (see Farmer, 2017).
- Basic problem reframed:

$$V(x_t, z_t) = \max_{x_{t+1} \in \Gamma(x_t, z_t)} F(x_t, z_t, x_{t+1}) + \beta \sum_{z_{t+1}} P(z_{t+1}|z_t) V(x_{t+1}, z_{t+1})$$

- Effects of discretization:
 - Choose x_{t+1} directly instead of y_t (can't leave grid).
 - Expectation is matrix multiplication.
- Notation: X(a, b; c) is a matrix where the columns stack over a and b (i.e., (a₁, b₁), (a₁, b₂), ..., (a₂, b₁), ...) and the rows stack over c.

Step 1: given iteration *k* guess **V**_{*k*}, optimize decision.

Define

$$\mathbf{Q}(x_t, z_t; x_{t+1}) = \underbrace{\mathbf{F}(x_t, z_t; x_{t+1})}_{NK \times N} + \beta \left(\underbrace{\mathbf{P}(z_t; z_{t+1})}_{K \times K} \otimes \underbrace{\mathbf{1}_N}_{N \times 1}\right) \underbrace{\mathbf{V}(x_{t+1}; z_{t+1})'}_{K \times N}$$

if $x_{t+1} \in \Gamma(x_t, z_t)$, and $-\infty$ otherwise.

 \blacktriangleright Reminder: \otimes is the Kronecker product, so that

$$A \otimes B = \begin{bmatrix} A_{11}B & \cdots & A_{1n}B \\ \vdots & \ddots & \vdots \\ A_{n1}B & \cdots & A_{nn}B \end{bmatrix}$$

• Define $x_{t+1}^*(x_t, z_t) = \arg \max_{x_{t+1}} \mathbf{Q}(x_t, z_t; x_{t+1})$. This is the column-wise max.

Step 2: given decisions, update V ("Howard Improvement").

- Can update \mathbf{V}_{k+1} by plugging in x_{t+1}^* and \mathbf{V}_k on the RHS, then iterate, but this is slow.
- Better approach: solve for **exact** value function under policy x_{t+1}^* .

Define:

$$\begin{aligned} \mathsf{A}(x_t, z_t, x_{t+1}, z_{t+1}) &= \mathsf{P}(z_{t+1}|z_t) \cdot \mathbf{1} \{ x_{t+1} = x_{t+1}^*(x_t, z_t) \} \\ \mathsf{F}^*(x_t, z_t) &= \mathsf{F}(x_t, z_t, x_{t+1}^*) \end{aligned}$$

Then we have:

$$\underbrace{\mathbf{V}(\mathbf{x}_{t}, \mathbf{z}_{t})}_{NK \times 1} = \underbrace{\mathbf{F}^{*}(\mathbf{x}_{t}, \mathbf{z}_{t})}_{NK \times 1} + \beta \underbrace{\mathbf{A}(\mathbf{x}_{t}, \mathbf{z}_{t}; \mathbf{x}_{t+1}, \mathbf{z}_{t+1})}_{NK \times NK} \underbrace{\mathbf{V}(\mathbf{x}_{t+1}, \mathbf{z}_{t+1})}_{NK \times 1}$$

which implies the exact solution $\mathbf{V}_{k+1} = (\mathbf{I} - \beta \mathbf{A})^{-1} \mathbf{F}^*$.

- lterate on Steps 1 and 2 until x_{t+1}^* stops changing. Then you are done!
- Stationary distribution: eigenvector of **A**' associated with unit eigenvalue.
 - Similarly, stationary dist. of exogenous states is eigenvector of P' with unit eigenvalue.
- Note: **A** will contain many zeros, often better to use sparse matrices.
- ▶ For *P*, use Rouwenhorst (1995) method to approximate Gaussian AR(1) processes.
 - Other approximations struggle as ho
 ightarrow 1.
 - Better to read treatment in Kopecky and Suen, RED 2010.
- Suffers from curse of dimensionality, but GPUs can provide huge speedup!

Special Case: Exogenous Asset Pricing

- Assume that all states are exogenous.
- Combine $E_t[M_{t+1}R_{t+1}] = 1$ and definition $R_{t+1} = (P_{t+1} + D_{t+1})/P_t$ to obtain

$$PD(z_t) = E_t \left\{ M(z_t, z_{t+1}) \left(PD(z_{t+1}) + 1 \right) \frac{D(z_{t+1})}{D(z_t)} \right\}.$$

Then we can solve for PD exactly with a single linear inversion:

$$\begin{split} A(z_t, z_{t+1}) &\equiv P(z_t, z_{t+1}) M(z_t, z_{t+1}) \frac{D(z_{t+1})}{D(z_t)} \\ PD(z_t) &= A(z_t; z_{t+1}) \left(PD(z_{t+1}) + \mathbf{1}_K \right) \\ PD &= (I - \mathbf{A})^{-1} \mathbf{1}_K \end{split}$$

Dynamic Programming: Continuous Models

Generic Optimality Conditions

- As long as the problem is well-behaved (uniquely determined by FOCs), it is usually better to solve the FOCs than to directly use the value function.
- > Typical approach is to just start taking derivatives, but can actually be more systematic.
- Let's add some additional structure (slight change of notation):
 - Let c_t be consumption, and y_t be all **other** controls.
 - Let $\Psi(x_t, c_t, y_t, z_t) \ge o$ be the budget constraint, and $\Gamma(x_t, c_t, y_t, z_t) \ge o$ be all other constraints.
 - Assume the budget constraint is written $c_t \leq \cdots$ so that $\partial \Psi_t / \partial c_t = -1$.
 - Let $F(x_t, y_t, z_t) = u(x_t, c_t, y_t, z_t)$.

Generic Optimality Conditions

Generic optimality condition for *y*_t:



- All quantities expressed in units of consumption.
- Marginal continuation values Ω_t defined by fixed point

$$\Omega_{t} = E_{t} \left\{ M_{t+1} \left[\left(\frac{\partial u_{t+1}}{\partial c_{t+1}} \right)^{-1} \frac{\partial u_{t+1}}{\partial x_{t+1}} + \frac{\partial \Psi_{t+1}}{\partial x_{t+1}} + \mu_{t+1} \frac{\partial \Gamma_{t+1}}{\partial y_{t+1}} + \Omega_{t+1} \frac{\partial x_{t+2}}{\partial x_{t+1}} \right] \right\}$$

where M_{t+1} is the SDF. Note: works for EZW preferences.

Example: Kaltenbrunner and Lochstoer (2010)

► Preferences:
$$U_t = \left((1-\beta)C_t^{1-\rho} + \beta E_t \left[U_{t+1}^{1-\gamma} \right]^{\frac{1-\rho}{1-\gamma}} \right)^{\frac{1}{1-\rho}}$$

► Budget constraint: $C_t \leq Z_t^{1-\alpha} K_t^{\alpha} - i_t K_t$.

- Endogenous state LOM: $K_{t+1} = (1 \delta)K_t + \phi(i_t)K_t$.
- Exogenous state LOM: $\log Z_{t+1} = \phi \log Z_t + \varepsilon_{t+1}$.
- Optimality conditions:

$$\mathbf{O} = -\mathbf{1} + \phi'(i_t)\Omega_{K,t}$$
$$\Omega_{K,t} = E_t \left\{ M_{t+1} \left[\alpha \left(\frac{Z_{t+1}}{K_{t+1}} \right)^{1-\alpha} - i_{t+1} + \left((1-\delta) + \phi(i_{t+1}) \right) \Omega_{K,t+1} \right] \right\}$$

Example: Kaltenbrunner and Lochstoer (2010)

• Preferences:
$$U_t = \left((1-\beta)C_t^{1-\rho} + \beta E_t \left[U_{t+1}^{1-\gamma} \right]^{\frac{1-\rho}{1-\gamma}} \right)^{\frac{1}{1-\rho}}$$

• Budget constraint: $C_t \leq Z_t^{1-\alpha} K_t^{\alpha} - i_t K_t$.

- Endogenous state LOM: $K_{t+1} = (1 \delta)K_t + \phi(i_t)K_t$.
- Exogenous state LOM: $\log Z_{t+1} = \phi \log Z_t + \varepsilon_{t+1}$.
- Optimality conditions:

$$1 = E_t [M_{t+1}R_{t+1}]$$

$$R_{t+1} \equiv \frac{\alpha (Z_{t+1}/K_{t+1})^{1-\alpha} - i_{t+1} + ((1-\delta) + \phi(i_{t+1}))q_{t+1}}{q_t}$$

$$q_t \equiv \phi'(i_t)^{-1}$$

Spring 2024

Complementary Slackness

• Complementary slackness: given constraint Γ_t and multiplier μ_t :

$$\mu_t \Gamma_t = 0, \qquad \mu_t \ge 0, \qquad \Gamma_t \ge 0.$$

- Example: lower bound $y_t \ge 0$.
 - Challenge: kinked, nondifferentiable policy function.
- Auxiliary variable (Garcia and Zangwill) approach:
 - Define policy function as auxiliary variable α_t .
 - Define $y_t = \max(\alpha_t, 0)^k$ for k > 1.
 - Define $\mu_t = \max(-\alpha_t, \mathbf{0})^k$ for k > 1.

Delivers continuously differentiable policy function.

Time Iteration

Assume equilibrium conditions follow $f(x, y, z, \mathcal{E}(x, y, z)) = 0$, where

$$\mathcal{E}_{t} = E_{t} \left[q(x_{t+1}, y_{t+1}, z_{t+1}) \right]$$

- Choose grid $\{\bar{x}_i, \bar{z}_i\}$ and basis functions $\psi(s, z)$.
- Let b_k be the coefficients from the previous (*k*th) iteration.
- Key idea: use previous guess b_k to form expectations \mathcal{E} :

$$y_{t+1} = b'_k \psi(x_{t+1}, z_{t+1})$$

$$\mathcal{E}_t = \sum_j \omega_j q(x_{t+1}, y_{t+1}, z_{t+1})$$

where ω_i are quadrature weights, then solve for y_t given \mathcal{E}_t .

Time Iteration on Controls

Time iteration on controls: for each (\bar{x}_i, \bar{z}_i) , choose y_i so $f(\bar{x}_i, y_i, \bar{z}_i, \mathcal{E}(\bar{x}_i, \bar{z}_i; b_k)) = 0$, with

$$\mathcal{E}_t(\bar{x}_i, \bar{z}_i; b_k) = \sum_j \omega_j q\Big(x_{t+1}, b'_k \psi(x_{t+1}, z_{t+1}), z_{t+1}\Big)$$

Recipe: given y_i , quadrature node $\bar{\varepsilon}_i$, compute

$$\begin{aligned} x_{t+1} &= g(\bar{x}_i, y_i, \bar{z}_i) \\ z_{t+1} &= h(\bar{z}_i, \bar{\varepsilon}_j). \end{aligned}$$

Use nonlinear equation solver. Gradient:

$$\frac{df}{dy_t} = \frac{\partial f}{\partial y_t} + \frac{\partial f}{\partial \mathcal{E}_t} E_t \left[\frac{\partial x_{t+1}}{\partial y_t} \left(\frac{\partial q}{\partial x_{t+1}} + \frac{\partial q}{\partial y_{t+1}} \frac{\partial y_{t+1}}{\partial x_{t+1}} \right) \right]$$

where $\partial y_{t+1} / \partial x_{t+1} = b'_k (\partial \psi(x_{t+1}, z_{t+1}) / \partial x_{t+1}).$

• Once solutions $\{y_i^*\}$ have been found, choose b_{k+1} so that $y_i^* = b'_{k+1}\psi(\bar{x}_i, \bar{z}_i) \ \forall i$, repeat.

Time Iteration

Iterating on controls is straightforward but not very efficient.

- Additional challenge: y_{t+1} may not be non-smooth in states.
- > Alternative: time iteration on coefficients. Solve directly for \hat{b} that satisfies

$$f\left(\bar{x}_{i},\hat{b}'\psi(\bar{x}_{i},\bar{z}_{i}),\bar{z}_{i},\mathcal{E}\right)=0.$$

i.e., compute $y_i = \hat{b}' \psi(\bar{x}_i, \bar{z}_i)$ and proceed as before.

Use nonlinear equation solver with gradient

$$\frac{\partial f}{\partial \hat{b}} = \left\{ \frac{\partial f}{\partial y_t} + \frac{\partial f}{\partial \mathcal{E}_t} E_t \left[\frac{\partial x_{t+1}}{\partial y_t} \left(\frac{\partial q}{\partial x_{t+1}} + \frac{\partial q}{\partial y_{t+1}} \frac{\partial y_{t+1}}{\partial x_{t+1}} \right) \right] \right\} \frac{\partial y_t}{\partial \hat{b}}$$

/ $\partial \hat{b} = \psi(x_t, z_t)'.$

• Update $b_{k+1} = \hat{b}$ and repeat until $||b_{k+1} - b_k||$ is smaller than some threshold.

where ∂v_t

Direct Solution

- Most efficient (but least robust): solve for b directly.
- Apply same solution to both sides:

$$O = f\left(x_t, b'\psi(x_t, z_t), z_t, \mathcal{E}_t\right)$$
$$\mathcal{E}_t = E_t\left[q\left(x_{t+1}, b'\psi(x_{t+1}, z_{t+1}), z_{t+1}\right)\right].$$

Run nonlinear equation solver with gradient

$$\frac{\partial f}{\partial b} = \frac{\partial f}{\partial y_t} \frac{\partial y_t}{\partial b} + \frac{\partial f}{\partial \mathcal{E}_t} E_t \left[\left(\frac{\partial q}{\partial x_{t+1}} + \frac{\partial q}{\partial y_{t+1}} \frac{\partial y_{t+1}}{\partial x_{t+1}} \right) \frac{\partial x_{t+1}}{\partial y_t} \frac{\partial y_t}{\partial b} + \frac{\partial q}{\partial y_{t+1}} \frac{\partial y_{t+1}}{\partial b} \right]$$

Can start with time iteration and then switch to direct solution to get best of both worlds.

Special Case: Exogenous Asset Pricing

Return to special case

$$PD(z_t) = E_t \left\{ M(z_t, z_{t+1}) \left(PD(z_{t+1}) + 1 \right) \frac{D(z_{t+1})}{D(z_t)} \right\}.$$

Apply guess $PD(z_t) = \psi(z_t)'b$, and use quadrature scheme $(\omega_j, \bar{\varepsilon}_j)$:

$$\psi(\mathbf{z}_t)'\mathbf{b} = \sum_j \omega_j \mathbf{M}(\mathbf{z}_t, \bar{\varepsilon}_j) \left(\frac{\mathbf{D}(\mathbf{z}_t, \bar{\varepsilon}_j)}{\mathbf{D}(\mathbf{z}_t)}\right) \left(\psi(\mathbf{z}_t, \bar{\varepsilon}_j)'\mathbf{b} + \mathbf{1}\right)$$

with slight abuse of notation to substitute $(z_t, \bar{\varepsilon}_j)$ for z_{t+1} .

Special Case: Exogenous Asset Pricing

If we now define

$$\begin{split} \mathsf{A}(\mathsf{z}_{\mathsf{t}}) &= \sum_{j} \omega_{j} \mathsf{M}(\mathsf{z}_{\mathsf{t}}, \bar{\varepsilon}_{j}) \left(\frac{\mathsf{D}(\mathsf{z}_{\mathsf{t}}, \bar{\varepsilon}_{j})}{\mathsf{D}(\mathsf{z}_{\mathsf{t}})} \right) \psi(\mathsf{z}_{\mathsf{t}}, \bar{\varepsilon}_{j}) \\ \mathsf{c}(\mathsf{z}_{\mathsf{t}}) &= \sum_{j} \omega_{j} \mathsf{M}(\mathsf{z}_{\mathsf{t}}, \bar{\varepsilon}_{j}) \left(\frac{\mathsf{D}(\mathsf{z}_{\mathsf{t}}, \bar{\varepsilon}_{j})}{\mathsf{D}(\mathsf{z}_{\mathsf{t}})} \right) \end{split}$$

then we obtain

$$\Psi b = \mathbf{A}b + \mathbf{C}$$

 $b = (\Psi - \mathbf{A})^{-1}\mathbf{C}$

Another one-step linear solution!

Additional Refinements

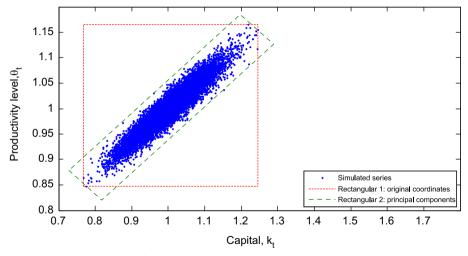
Precomputation:

- Can save time by pre-computing $\psi(\bar{x}_i, \bar{z}_i)$.
- Note that for a given grid $\{\bar{z}_i\}$ and nodes $\{\bar{z}_i\}$, end up with the same grid over z_{t+1} .
- If $\psi(\mathbf{x}_{t+1}, \mathbf{z}_{t+1}) = \psi_{\mathbf{x}}(\mathbf{x}_{t+1})\psi_{\mathbf{z}}(\mathbf{z}_{t+1})$, then we can also precompute $\psi_{\mathbf{z}}(\mathbf{z}_{t+1}(\bar{\mathbf{z}}_{i}, \bar{\mathbf{z}}_{j}))$.

Adaptive domain:

- Approximations work better when variables are not highly correlated.
- Better: use principal components as states.
- Do SVD to obtain $X = (x_1, \ldots, x_T)' = USV'$, then the PCs are $\tilde{X} = XV$.
- Recover X using $X = \tilde{X}V'$.

Adaptive Domain: Illustration



Source: Judd, Maliar, Maliar, Valero (2013).

Special Case: Endogenous Grid Method

> Typically choose grid, then solve so optimality conditions hold on it.

- But sometimes can skip optimization step by exploiting properties of equilibrium condition.
- Endogenous grid method of Carroll (2006).

Example: consider a life-cycle Bewley model with Euler equation

$$c_t(a_t, y_t)^{-\gamma} = \beta E_t \left[c_{t+1}(a_{t+1}, y_{t+1})^{-\gamma} \right]$$

where t is age, a is assets, and y is income, subject to the budget constraint

$$c_t + R^{-1}a_{t+1} = a_t + y_t$$

• Given (a_t, y_t) , we cannot solve for c_t (equivalently, a_{t+1}) in closed form.

Special Case: Endogenous Grid Method

But what if we somehow knew next period's value of a_{t+1} and next period's policy function c_{t+1} ? Then from the Euler equation we would know

$$c_{t}^{*} = \left\{\beta E_{t} \left[c_{t+1} (a_{t+1}, y_{t+1})^{-\gamma} \right] \right\}^{-1/\gamma}$$

and from the budget constraint we would know

$$a_t^* = c_t^* + R^{-1}a_{t+1} - y_t.$$

- This means that if we start at (a_t^*, y_t) , c_t^* is the optimal policy!
- For a given grid of a_{t+1} values, we can solve for (c_t^*, a_t^*, y_t) , then approximate $c_t(a_t, y_t)$ on this grid (typically by linearly interpolating).
- > Not a generic method, but when it works it is very simple and effective.
 - See Maliar and Maliar (2013), and other work for similar envelope condition method.

Perturbation Methods

Perturbation Methods

- Based on local expansion around a point.
- Computationally cheap, but less accurate far from approximation point.
 - Great as initial guess for global solution.
- Fold exogenous states into *x*_t to rewrite

$$\mathbf{O} = E_t \left[f(\mathbf{x}_t, \mathbf{y}_t, \mathbf{x}_{t+1}, \mathbf{y}_{t+1}) \right]$$

$$\mathbf{y}_t = g(\mathbf{x}_t, \sigma)$$

$$\mathbf{x}_{t+1} = h(\mathbf{x}_t, \sigma) + \sigma \eta \varepsilon_{t+1}.$$

Note that g and h are different from earlier notation.

Perturbation Methods

First order perturbation:

$$y_t = g(\bar{x}, \mathbf{O}) + g_x(\bar{x}, \mathbf{O})(x - \bar{x}) + g_\sigma(\bar{x}, \mathbf{O})\sigma$$

$$x_{t+1} = h(\bar{x}, \mathbf{O}) + h_x(\bar{x}, \mathbf{O})(x - \bar{x}) + h_\sigma(\bar{x}, \mathbf{O})\sigma + \sigma\eta\varepsilon_{t+1}$$

Choose values to set derivatives of equilibrium condition to zero:

$$\begin{aligned} &(\mathbf{x}_t): \quad \mathbf{o} = \mathbf{f}_{\mathbf{x}} + \mathbf{f}_{\mathbf{y}} \mathbf{g}_{\mathbf{x}} + \mathbf{f}_{\mathbf{x}'} \mathbf{h}_{\mathbf{x}} + \mathbf{f}_{\mathbf{y}'} \mathbf{g}_{\mathbf{x}} \mathbf{h}_{\mathbf{x}} & (\mathbf{1}) \\ &(\sigma): \quad \mathbf{o} = \mathbf{f}_{\mathbf{y}} \mathbf{g}_{\sigma} + \mathbf{f}_{\mathbf{x}'} \mathbf{h}_{\sigma} + \mathbf{f}_{\mathbf{y}'} \left(\mathbf{g}_{\sigma} + \mathbf{g}_{\mathbf{x}} \mathbf{h}_{\sigma} \right) & (\mathbf{2}) \end{aligned}$$

- Solution to (2) implies $g_{\sigma} = h_{\sigma} = 0$ (no risk effects).
- Apply $g(\bar{x}, o) = \bar{y}$, $h(\bar{x}, o) = \bar{x}$, define e.g., $\hat{x} = x \bar{x}$, to obtain system that must solve (1):

$$\hat{y}_t = g_x \hat{x}_t$$
$$\hat{x}_{t+1} = h_x \hat{x}_t + \sigma \eta \varepsilon_{t+1}$$

Many solution techniques: Sims (2001), Klein (2000).

Higher Order Perturbations

Second-order perturbation (see e.g., Judd and Guu (1997) for solution method):

$$\begin{split} \hat{y}_t &= \mathbf{g}_{\mathbf{x}} \hat{x}_t + \frac{1}{2} \mathbf{G}_{\mathbf{x}\mathbf{x}} (\hat{x}_t \otimes \hat{x}_t) + \frac{1}{2} \mathbf{g}_{\sigma\sigma} \sigma^2 \\ \hat{x}_{t+1} &= \mathbf{h}_{\mathbf{x}} \hat{x}_t + \frac{1}{2} \mathbf{H}_{\mathbf{x}\mathbf{x}} (\hat{x}_t \otimes \hat{x}_t) + \frac{1}{2} \mathbf{h}_{\sigma\sigma} \sigma^2 + \sigma \eta \varepsilon_{t+1} \end{split}$$

- Now risk influences policy functions (in a constant way).
 - Third-order: $\sigma^2 \hat{x}_t$ term linear in states.
 - Higher order: nonlinear risk-state interactions.
- Major problem: explosiveness. Univariate example:

$$\hat{x}_{t+1} = \dots + \frac{1}{2}h_{xx}\hat{x}_t^2 = \dots + \frac{1}{2}h_{xx}\left(\dots + \frac{1}{2}h_{xx}\hat{x}_{t-1}^2\right)^2$$

Pruned State Space

Pruned state-space approach (Andreasen et al, 2018). Split xt into first-order terms xt and second-order terms xt:

$$\begin{split} \hat{\mathbf{x}}_{t+1}^{f} &= \mathbf{h}_{\mathbf{x}} \hat{\mathbf{x}}_{t}^{f} + \sigma \eta \varepsilon_{t+1} \\ \hat{\mathbf{x}}_{t+1}^{s} &= \mathbf{h}_{\mathbf{x}} \hat{\mathbf{x}}_{t}^{s} + \frac{1}{2} \mathbf{H}_{\mathbf{X}\mathbf{X}} (\hat{\mathbf{x}}_{t}^{f} \otimes \hat{\mathbf{x}}_{t}^{f}) + \frac{1}{2} \mathbf{h}_{\sigma\sigma} \sigma^{2} \end{split}$$

- ▶ No interaction between x^s and x^s means no explosiveness: $\hat{x}_t = A(L)\varepsilon_t + B(L)\varepsilon_t^2$.
- Policy functions:

$$\hat{y}_t = \mathbf{g}_{\mathbf{x}} \left(\hat{x}_t^f + \hat{x}_t^s \right) + \frac{1}{2} \mathbf{G}_{\mathbf{x}\mathbf{x}} (\hat{x}_t^f \otimes \hat{x}_t^f) + \frac{1}{2} \mathbf{g}_{\sigma\sigma} \sigma^2$$

See paper for third-order equivalent.

- Perfect foresight paths (also known as deterministic transition paths, or "MIT shocks")
- Idea: if we assume no risk from today on, then path back to steady state is solution to equilibrium conditions.
- Notation for equilibrium conditions:

 $f(s_{t-1}, s_t, s_{t+1}; z_t) = 0$

where s' = (x', y') are endogenous states and policy functions, and z are exogenous states.

- > Deterministic environment buys a lot of tractability.
 - Because no shocks will arrive, we can directly use s_{t+1} . Don't need expectations.
 - Can directly use nonlinear equilibrium conditions for *f*, no need to linearize.
 - Can change parameters or apply shocks to exogenous states.

- Solution and notation follow Juillard, Laxton, McAdam, Pioro (1998).
- Stack equations to form

$$\mathbf{f}(\mathbf{s}) = \begin{bmatrix} f_{0}(s_{0}) \\ f_{1}(s_{0}, s_{1}, s_{2}) \\ \vdots \\ f_{T}(s_{T-1}, s_{T}, s_{T+1}) \\ f_{T+1}(s_{T+1}) \end{bmatrix} = \mathbf{0}$$

including additional initial and terminal equations, typically

$$f_0(s_0) = s_0 - s_0^*$$
 $f_{T+1}(s_{T+1}) = s_{T+1} - s_{T+1}^*$

where s_0^* is the initial steady state, and s_{T+1}^* is the final steady state.

(3)

Typically solved using Newton's method:

$$\underbrace{\Delta \mathbf{s}}_{\text{step size}} = -\mathbf{f}'(\mathbf{s})^{-1}\mathbf{f}$$

• Key to solution is computing inverse Jacobian $\mathbf{f}'(\mathbf{s})^{-1}$. Want to solve:

$$\begin{bmatrix} I & & & \\ L_1 & C_1 & F_1 & & \\ & \ddots & \ddots & \ddots & \\ & & L_T & C_T & F_T \\ & & & & I \end{bmatrix} \Delta \mathbf{s} = -\mathbf{f}$$

where L_t , C_t , and F_t are the derivatives of $f_t(s_{t-1}, s_t, s_{t+1})$ with respect to s_{t-1} , s_t , s_{t+1} .

- ▶ In practice, this matrix has size $n(T+2) \times n(T+2)$, where *n* is the number of equilibrium conditions and *T* is the number of periods.
 - Most of the entries are zeros, so sparse matrix tools can handle it.
 - Alternative: Juillard, Laxton, McAdam, Pioro (1998) provide a recursive algorithm computing Δs .
- Weakness of the approach: exactly end at steady state.
 - May require huge number of periods to avoid distorting the calculations.
- > My alternative: assume that by end of the sample equilibrium follows linearized solution.
 - Linearized solution: $y_t = G_x x_t + G_z z_t$ where y_t are endog. controls and x_t are endog. states.
 - Replace terminal condition with the following (*h* is transition equation):

$$f_{T+1}(s_t) = \begin{bmatrix} x_{T+1} - h(s_T, z_{T+1}) \\ y_{T+1} - G_x x_{T+1} - G_z z_{T+1} \end{bmatrix} = 0.$$

- Some questions require heterogeneous agent models.
 - Although you should keep in mind that some do not.
- In these cases, working with endogenous aggregate states is complex.
 - Often, only a small subset of aggregate quantities (e.g., prices) matter for individual behavior.
 - However, values of these aggregates may depend on the entire distribution.
 - Krusell-Smith approach approximates using simpler forecasts based only on moments.
 - But computationally intensive, and no guarantee this will work well.
- Recent alternative: sequence space jacobian.
 - Method to compute linearized impulse responses or perfect foresight paths.
 - Note: these solutions remove aggregate risk, but not idiosyncratic risk.

- Start with a function that defines the aggregate equilibrium H = 0.
- Example in neoclassical production model, capital market clearing:

$$H_t(K,Z) = \int_i k_{i,t}(X,Z) - K_t.$$

► For linearized impulse response, can use approximation

 $H_K dK + H_Z dZ = 0$

to obtain

$$dK = -H_K^{-1}H_Z dZ$$

- Can also solve this H(K, Z) = 0 as nonlinear system of equations.
- Key to both solutions is the Jacobian, (H_K, H_Z) .

- First, we need to split the problem.
 - Het. agent models generally intractable when behavior depends on entire distribution.
 - Need to collapse to a subset of aggregate states X_t sufficient for the individual's problem.
 - In classical Krusell-Smith model, this is just prices (r_t, w_t) .

b Define **block** Y = h(X) to be mapping between sufficient states X and aggregate outputs Y.

- In this example, X is prices, Y is capital demand.
- Full model equilibrium requires multiple blocks:

$$H(K,Z) = H(h(X),Z) = H(h(g(K,Z))Z)$$

where g(K, Z) maps states into prices (equal marginal products from firm FOCs).

- Efficiency gains from applying closed form solutions when available (see paper).
- ▶ Then can compute Jacobian (H_K, H_Z) using the chain rule given Jacobians of $h, X(\cdot)$.

Define notation for the problem as

Individual optimality: Distribution law of motion: Measurement of agg. states:

$$\begin{split} \mathbf{v}_t &= \mathbf{v}(\mathbf{v}_{t+1}, \mathbf{X}_t) \\ \mathbf{D}_{t+1} &= \Lambda(\mathbf{v}_{t+1}, \mathbf{X}_t)' \mathbf{D}_t \\ \mathbf{Y}_t &= y(\mathbf{v}_{t+1}, \mathbf{X}_t)' \mathbf{D}_t. \end{split}$$

- Apply a single shock of size dx to X at time s.
 - Then we want to compute dY_t^s , change in Y at time t due to shock at time s.
- ► Take limit as $dx \rightarrow 0$: $dY_t^s = (dy_t^s)' D_t^s + (y_t^s)' dD_t^s$
- Possible (but costly) to compute directly.
 - Apply the shock at time *s*, solve **v** backwards, then iterate **D** forwards.
 - Repeat this for each time s.

First efficiency gain: policy functions depend only on distance to shock s - t

$$\mathbf{y}_t^{\mathsf{s}} = \mathbf{y}_{t+k}^{\mathsf{s}+k}, \qquad \qquad \Lambda_t^{\mathsf{s}} = \Lambda_{t+k}^{\mathsf{s}+k}.$$

Second gain: use the fact that $dx \rightarrow o$ to simplify the problem

$$dY_t^s = (d\mathbf{y}_t^s)'\mathbf{D}_t^s + (\mathbf{y}_t^s)'d\mathbf{D}_t^s = (d\mathbf{y}_t^s)'\mathbf{D}_{ss} + \mathbf{y}_{ss}'d\mathbf{D}_t^s.$$

Now subtract dY_{t-1}^{s-1} :

$$dY_t^{s} - dY_{t-1}^{s-1} = \underbrace{(d\mathbf{y}_t^{s} - d\mathbf{y}_{t-1}^{s-1})'}_{=o} \mathbf{D}_{ss} + \mathbf{y}_{ss}'(d\mathbf{D}_t^{s} - d\mathbf{D}_{t-1}^{s-1})$$
$$= \mathbf{y}_{ss}'(d\mathbf{D}_t^{s} - d\mathbf{D}_{t-1}^{s-1})$$

Spring 2024

Difference in distributions:

$$d\mathbf{D}_{t}^{s} = (d\Lambda_{t-1}^{s})'\mathbf{D}_{ss} + (\Lambda_{ss})'d\mathbf{D}_{t-1}$$

Now difference as in previous slide:

$$d\mathbf{D}_{t}^{s} - d\mathbf{D}_{t-1}^{s-1} = \underbrace{(d\Lambda_{t-1}^{s} - d\Lambda_{t-2}^{s-1})'}_{=o} \mathbf{D}_{ss} + \Lambda_{ss}'(d\mathbf{D}_{t-1} - d\mathbf{D}_{t-2})$$
$$= \Lambda_{ss}'(d\mathbf{D}_{t-1} - d\mathbf{D}_{t-2})$$
$$\vdots$$
$$= (\Lambda_{ss}')^{t-1}(d\mathbf{D}_{1} - d\mathbf{D}_{o})$$
$$= (\Lambda_{ss}')^{t-1}(d\Lambda_{0}^{1})'\mathbf{D}_{ss}$$

Recursive structure re-using repeated terms much cheaper to compute.

Conclusion

Many tools available, want to select right tools for the right job.

- More complex or high tech is not always better!
 - Simpler models are often easier to understand.
 - You can run lots of things to gain intuition about role of different mechanisms.
 - You retain degrees of freedom to use on other features.
- > My advice: start with simple methods before complexifying.
 - My personal favorite: perfect foresight paths.