SOLVING DYNAMIC MODELS

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Welcome!

- ▶ Why solve dynamic models?
- ► Theory vs practice
- Resources

The basic dynamic programming setup

A firm's dynamic price adjustment problem

Value function iteration

Discretization

Simulation

A DYNAMIC OPTIMIZATION PROBLEM

$$\max_{\{s_{t+k+1}\}_{k=0}^{\infty}} \mathbb{E}_t \sum_{k=0}^{\infty} \beta^k R(s_{t+k}, x_{t+k}, s_{t+k+1})$$
s.t. $s_{t+k+1} \in \Gamma(s_{t+k}, x_{t+k}) \quad \forall k$

$$x_{t+k} | x_{t+k-1}, \dots \sim F(x_{t+k} | x_{t+k-1})$$

- Decisionmaker maximizes an expected present discounted value with discount rate $0 < \beta < 1$ by choosing a contingent plan.
- ▶ Current payoff $R(\cdot)$ in t depends on potentially vector-valued:
 - ightharpoonup State s_t
 - ightharpoonup Stationary Markov shock x_t
 - ightharpoonup *Policy* s_{t+1} , i.e., future state
- Policy choice is subject to constraints $\Gamma(\cdot)$, which usually embed a "today vs tomorrow" dynamic tradeoff

Note: Lots of stuff fits into this structure. Household savings for welfare maximization, firm investment, financing, etc...

THE KEY RECURSIVE INSIGHT OF DYNAMIC PROGRAMMING

Sequence Formulation

$$\max_{\{s_{t+k+1}\}_{k=0}^{\infty}} \mathbb{E}_t \sum_{k=0}^{\infty} \beta^k R(s_{t+k}, x_{t+k}, s_{t+k+1}) =$$

$$\underbrace{\max_{\left\{s_{t+k+1}\right\}_{k=0}^{\infty}} R(s_{t}, x_{t}, s_{t+1}) + \beta \mathbb{E}_{t} \left[\underbrace{R(s_{t+1}, x_{t+1}, s_{t+2}) + \beta \mathbb{E}_{t+1} R(s_{t+2}, x_{t+2}, s_{t+3}) + \dots}_{V(s_{t+1}, x_{t+1})}\right]}_{V(s_{t}, x_{t})}$$

Stationary Bellman Equation, Fixed Point, or Recursive Formulation

$$V(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'|x} V(s',x') \right]$$
$$V = \mathcal{T}(V), \quad \mathcal{T}(f) = \max[R + \beta \mathbb{E}(f)]$$

THE GOAL

Starting with a Bellman equation formulation

$$V(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'|x} V(s',x') \right],$$

the goal is to introduce computational methods allowing us to

- **solve** the model for the optimal policy function s'(s,x), which usually involves solving for the value function V(s,x), and
- **simulate** the model by drawing exogenous Markov shock processes $x_1, x_2,$ which lead to endogenous sequences of states $s_1, s_2,$

Note: The goals aren't picked randomly. Structural estimation embeds them within each evaluation of an objective function.

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A CONCRETE EXAMPLE

A firm facing downward-sloping demand and marginal cost shocks chooses prices to maximize the expected PDV of its payouts, which are variable profits net of price adjustment costs.

$$V(p_{-1}, m) = \max_{p} \left[\Pi(p, m) - AC(p_{-1}, p) + \beta \mathbb{E}_{m'|m} V(p, m') \right]$$
$$\log m' = \rho \log m + \sigma \eta', \quad \eta' \sim N(0, 1)$$
$$\Pi(p, m) = (p - m)p^{-\varepsilon}, \quad AC(p_{-1}, p) = \frac{c}{2}(p - p_{-1})^2$$
$$0 < \beta < 1, \quad 0 < \rho < 1, \quad \sigma > 0, \quad \varepsilon > 1, \quad c > 0$$

Stepping Back

- Why is this a dynamic problem rather than a series of identically structured static monopolistic pricing decisions?
- ▶ What do we expect intuitively from the shape and slopes of the solutions $p(p_{-1}, m)$ and $V(p_{-1}, m)$?

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SOLVING THE MODEL WITH VFI

Solving the Bellman equation

$$V(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'|x} V(s',x') \right]$$

is equivalent to finding a fixed point of the operator ${\mathcal T}$

$$\iff V = \mathcal{T}(V), \quad \mathcal{T}(f) = \max [R + \mathbb{E}(f)].$$

Under regularity conditions, a famous contraction mapping theorem guarantees that value function iteration (VFI) will converge, i.e., repeatedly iterating on $\mathcal{T}(\cdot)$ converges

$$V_{(1)}, V_{(2)} = \mathcal{T}(V_{(1)}), \dots, V_{(n)} = \mathcal{T}(V_{(n-1)}), \dots \to_{n \to \infty} V$$

Some Notes

- ▶ I'm skipping over formalism. What regularity conditions? What is convergence in a function space? Does *V* exist uniquely? Why exactly are the sequence and Bellman formulations equivalent?...
- In each VFI step you compute optimal policies s'(s,x). So both values V(s,x) and optimal policies are obtained, as desired.

A BIT MORE DETAIL ON VFI

Start with R(s,x,s'), F(x'|x), $\Gamma(s,x)$, and $0<\beta<1$, as well as a guess $V_{(1)}$ and a specified solution tolerance $\varepsilon_{tol}>0$. For each step of VFI n=1,2,..., do

1. Compute $V_{(n+1)}$ by solving for optimal s' at each (s,x) via

$$V_{(n+1)}(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'\mid x} V_{(n)}(s',x') \right].$$

2. Compute the error

$$||V_{(n+1)} - V_{(n)}|| = \max_{s,x} |V_{(n+1)}(s,x) - V_{(n)}(s,x)|$$

3. If $||V_{(n+1)} - V_{(n)}|| < \varepsilon_{tol}$, exit. If $||V_{(n+1)} - V_{(n)}|| \ge \varepsilon_{tol}$, go to next iteration by jumping back to Step ?? with n = n + 1.

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Some of This Stuff is Tricky

Solving the general version of the Bellman equation

$$V(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'|x} V(s',x') \right]$$

involves some stuff that can be computationally tricky:

- **Functions** of continuous inputs like $V(\cdot)$ take some care to store, requiring interpolation methods, attention to smoothness, etc...
- **Expectations** of continuous Markov chains x'|x involve numerical integration or quadrature which can be tricky or time-consuming.
- **Optimization** over continuous choices s' requires application of potentially finicky and time-consuming numerical algorithms.

But if we **discretize** everything by putting s, x, and s' on grids, then

functions are easy vectors, expectations are easy weighted sums, and optimization is just picking the biggest number from a vector.

DISCRETIZING MARKOV CHAINS

We desire to work with methods converting from a general continuous, stationary Markov chain described by the distribution $F(x^\prime|x)$ to

- ▶ A grid $\{\bar{x}_1,...,\bar{x}_{N_x}\}$ of N_x points in which discretized x lives.
- ▶ A transition matrix Π^x with size $N_x \times N_x$ with entries

$$\Pi_{i,j}^x = \pi_{ij}^x = \mathbb{P}(x' = \bar{x}_j | x = \bar{x}_i)$$

Once we do this, note that integration collapses to a simple weighted sum:

$$\mathbb{E}_{F}\left[g(x')|x\right] = \int g(x')dF(x'|x) = \int g(x')f(x'|x)dx$$

$$\iff$$

$$\mathbb{E}_{\Pi^{x}}\left[g(x')|x = \bar{x}_{i}\right] \approx \sum_{j=1}^{N_{x}} \pi_{ij}^{x}g(\bar{x}_{j})$$

TAUCHEN (1986)

One classic simple method for discretizing the stationary AR(1) processes

$$x' = \rho x + \sigma \eta', \quad \eta' \sim N(0, 1), \quad 0 < \rho < 1, \quad \sigma > 0$$

comes from a famous paper Tauchen (1986).

- 1. Choose an (odd) value for N_x and a multiple M_x of standard deviations $\sigma_x = \frac{\sigma}{\sqrt{1-\rho^2}}$ which you want the grid to span.
- 2. Linearly space the grids points \bar{x}_i with gaps $\Delta > 0$

$$\{-M_x\sigma_x, -M_x\sigma_x + \Delta, ..., -\Delta, 0, \Delta, ..., M_x\sigma_x - \Delta, M_x\sigma_x\}$$

3. Compute the entries of Π^x by integrating under the error term, i.e.,

$$\Pi_{ij}^{x} = \mathbb{P}(x' = \bar{x}_j | x = \bar{x}_i) = H\left(\bar{x}_j + \frac{\Delta}{2}; \rho \bar{x}_i, \sigma^2\right) - H\left(\bar{x}_j - \frac{\Delta}{2}; \rho \bar{x}_i, \sigma^2\right)$$

where $H(\cdot;\mu,\sigma^2)$ is the $N(\mu,\sigma^2)$ CDF. Endpoints absorb the tails.

Note: Within Tauchen's method, you can add constants, make this a lognormal process, apply more complicated grids, do this with vector processes, etc... Also, many other fancier non-Tauchen methods now exist.

DISCRETIZED BELLMAN EQUATION

$$V(s,x) = \max_{s' \in \Gamma(s,x)} \left[R(s,x,s') + \beta \mathbb{E}_{x'|x} V(s',x') \right].$$

- ▶ Discretize the Markov chain x'|x with N_x points.
- ▶ Choose a grid of N_s points on which s lives $\{\bar{s}_1,...,\bar{s}_{N_s}\}$.
- ▶ The discretized Bellman equation for each (\bar{s}_i, \bar{x}_j) is

$$V(\bar{s}_{i}, \bar{x}_{j}) = \max_{\substack{s' \in \Gamma(\bar{s}_{i}, \bar{x}_{j}), \\ s' \in \{\bar{s}_{1}, ..., \bar{s}_{N_{s}}\}}} \left[R(\bar{s}_{i}, \bar{x}_{j}, s') + \beta \sum_{k=1}^{N_{x}} \pi_{j,k}^{x} V(s', \bar{x}_{k}) \right].$$

lacktriangle Collecting across all (i,j), we have the convenient vectorized form

$$\underbrace{ \mathbf{V}}_{N_s N_x \times 1 \text{vector}} = \underbrace{ \mathbf{RHS}}_{\text{row max of specially constructed } N_s N_x \times N_s \text{matrix} }$$

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SIMULATING YOUR MODEL

After model solution, we often want to examine simulated data.

- ▶ Simulate the exogenous Markov chain $\{x_t\}_{t=1}^T$ for a large T.
- ▶ Using optimal policies $s'(\cdot)$, compute the implied states $\{s_t\}_{t=1}^T$.
- Process the data $\{s_t, x_t\}_{t=1}^T$ in whichever manner is needed for your question of interest.

Some Notes

In practice, when simulating for application of structural estimation techniques you may need to:

- ▶ Compare U(0,1) shocks with thresholds implied by Π^x .
- Simulate data with a panel structure, i.e., (s_{it}, x_{it}) for firms i = 1, ..., N and t = 1, ..., T.
- ▶ Be careful with initial conditions, which can be overly influential for persistent processes, by discarding "burn-in" periods.
- Be careful to set seeds for replicability of random draws in the simulation of exogenous processes.

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