

Advanced Macroeconomics I

Chapter II: *Technical Preliminaries*

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Outline

1. Difference equations
2. Dynamic optimization

Introduction

- In this chapter, we provide an introduction into the main tools and techniques that will be applied throughout this course:
 - (Linear) difference equations
 - Linear state space models
 - (Stochastic) dynamic optimization techniques
 - Markov processes

Outline

1. Deterministic linear difference equations

1.1 Setup

1.2 Characterization

1.3 Solving LDEs

1.4 Coupled systems of higher order

1.5 Decoupling the system

2. Stochastic linear difference equations

2.1 Setup and stochastic structure

2.2 Conditional moments/forecasting

2.3 Unconditional moments and covariance stationarity

2.4 State-space systems and impulse response functions

First-order difference equations

- Consider the **first-order linear difference equation** (LDE)

$$x_t = ax_{t-1} + e_t, \quad t = 1, 2, \dots \quad (1)$$

where

- a is a constant coefficient/parameter
- x_t and e_t are scalar real-valued functions of time $t = 1, 2, \dots$:
 - $x_t = x(t)$
 - $e_t = e(t)$
- We assume **perfect foresight**, i.e. e_t is known for all t
- In economic applications, we will refer to x_t as the **endogenous variable** and e_t as the **exogenous forcing variable**
- e_t is assumed to be a bounded scalar real-valued sequence (generalization to vector difference equations straightforward)
- A **solution** to (1) is an expression giving x_t as a function only of past, present, and future values of e_t (and **boundary conditions**)

Boundedness

Definition 1 (Boundedness)

A sequence $\{X_t\}$ is bounded/stable if there exists $M > 0$ such that $\|X_t\|_\infty < M \forall t$

- Typically, only bounded solutions are economically relevant as otherwise, there is no scarcity
- Some problems, particularly those involving growth, may require appropriate stationarizing, e.g. transforming into intensive form

Characterizing difference equations: two important distinctions

- The stability of the LDE determines how to actually solve the LDE
 - Is the difference equation stable, i.e. $|a| \leq 1$?
→ solve the equation backward
 - Or do we have an unstable root, i.e. $|a| > 1$?
→ solve the equation forward
- The type of available boundary condition determines how we *should* solve the LDE:
 - Do we have an initial condition that restricts x to a particular value today or in the past?
→ solve the equation backward
 - Do we have a terminal condition that restricts x to a particular value in the future?
→ solve the equation forward
- Sometimes the type of stability condition does not match the available boundary condition: leads to problems with **uniqueness and existence of bounded solutions**

The lead/lag operator

- LDEs can generally be solved by **recursive substitution**, i.e. iteratively plugging in
- This can be made easier by introducing the lag/lead-operator L defined through

$$L^z ax_t = ax_{t-z}, \quad (2)$$

where a is a constant and z is an integer

- We will first consider the case of a stable difference equation, i.e. $|a| < 1$
- *Exercise 1*: Show that for a scalar $|a| < 1$ it holds that

$$\sum_{i=0}^{\infty} a^i L^i = (1 - aL)^{-1} = 1 + aL + a^2 L^2 + \dots \quad (3)$$

Solving backward

- To find a solution for x_t , rewrite (1) using the lag operator as

$$x_t(1 - aL) = e_t \quad (4)$$

- Exercise 2:* Show that the **non-homogenous component**, which expresses x_t as a function of current and past e_t

$$x_t = \sum_{s=-\infty}^t a^{t-s} e_s \quad (5)$$

is a solution

- Adding the **homogenous component** ca^t , with c being an arbitrary constant, delivers the general solution:

$$x_t = \sum_{s=-\infty}^t a^{t-s} e_s + ca^t = (1 - aL)^{-1} e_t + ca^t \quad (6)$$

- Exercise 3:* Show that this is an actual solution to the LDE
- Due to discounting with $|a| < 1$, this backward solution is bounded and the effect of past e_t decays geometrically

Ruling out indeterminacy using a boundary condition

- Problem: (6) will hold for any arbitrary c and there is no way to determine it without further information
- This case is called **indeterminacy**
- But: suppose we have an initial condition stating that x_0 is some given (known) constant.
- From the general solution (6) at time 0 we know that

$$x_0 = \sum_{s=-\infty}^0 a^{-s} e_s + ca^0, \quad (7)$$

- This equation then determines c uniquely as

$$c = x_0 - \sum_{s=-\infty}^0 a^{-s} e_s \quad (8)$$

The unique and determinate backward solution

- Inserting expression (8) for c into the solution for x_t , (6), we get

$$x_t = \sum_{s=-\infty}^t a^{t-s} e_s + ca^t \stackrel{(8)}{=} \sum_{s=-\infty}^t a^{t-s} e_s + a^t \left(x_0 - \sum_{s=-\infty}^0 a^{-s} e_s \right) \quad (9)$$

- Using that

$$\sum_{s=-\infty}^t a^{t-s} e_s - \sum_{s=-\infty}^0 a^{t-s} e_s = \sum_{s=1}^t a^{t-s} e_s \quad (10)$$

yields

$$x_t = \sum_{s=1}^t a^{t-s} e_s + a^t x_0 \quad (11)$$

- Due to discounting with $|a| < 1$, the backward solution with starting value x_0 is unique and bounded

Solving forward

- But what if $|a| > 1$, in which case the backward solution (11) will be unbounded?
- Fortunately, we can then try to solve forward
- Consider the difference equation (1) at time $t + 1$

$$x_{t+1} = ax_t + e_{t+1} \quad (12)$$

- *Exercise 4:* Show, using the lead operator, that the non-homogenous component of the forward solution is given by

$$x_t = - \sum_{s=t+1}^{\infty} \left(\frac{1}{a}\right)^{s-t} e_s \quad (13)$$

- Again, there exist more solutions that include the homogenous component ca^t :

$$x_t = - \sum_{s=t+1}^{\infty} \left(\frac{1}{a}\right)^{s-t} e_s + ca^t \quad (14)$$

- Since $|a| > 1$, unless $c = 0$, $ca^t \rightarrow \pm\infty$ for $t \rightarrow \infty$

Forward solutions and boundedness

- (14) doesn't look like much progress: we still cannot pin down c and the sequence for x_t grows without bounds for $t \rightarrow \infty$
- An initial condition as before would make the problem determinate, but still unbounded
- However: in economic applications sensible solutions typically require economic quantities to be bounded
- This amounts to a **terminal condition** $x_t < \infty$ for $t \rightarrow \infty$
- This leaves $c = 0$ as the only possible choice
→ the initial value x_0 of the sequence has to adjust to ensure that $c = 0$ is satisfied
- Sidenote: justification of terminal conditions can be subject of intense debate (e.g. Cochrane 2011)
- Physical capital stock obviously cannot go to infinity, but what about inflation?

Economic example: stable difference equation

- **Law of motion for capital** accumulation with given initial value k_0

$$k_{t+1} = (1 - \delta)k_t + i_{t+1} \quad (15)$$

- With $(1 - \delta) \in [0, 1)$, one can apply the backward solution (11):

$$\begin{aligned} k_t &= \sum_{s=-\infty}^t (1 - \delta)^{t-s} i_s + c(1 - \delta)^t \\ &= \sum_{s=1}^t (1 - \delta)^{t-s} i_s + (1 - \delta)^t k_0 \end{aligned} \quad (16)$$

- Capital stock at each point in time is the sum of
 1. what depreciation has left from the initial capital stock at $t = 0$
 2. the cumulative historical additions through (depreciated) investment
- National accounting often relies on this **perpetual inventory method**

Economic example: unstable difference equation

- No-arbitrage asset pricing relationship between
 - bond with constant coupon/interest rate $r > 0$ and
 - stock with purchase price p_t and known dividend d_t
- Interest rate on bond needs to equal dividend yield plus capital gains yield:

$$1 + r = \frac{d_{t+1} + p_{t+1}}{p_t} \quad (17)$$

- Since $(1 + r) > 1$, the resulting difference equation

$$p_{t+1} = (1 + r)p_t - d_{t+1} \quad (18)$$

can be solved forward using (14):

$$p_t = \sum_{s=t+1}^{\infty} \left(\frac{1}{1+r} \right)^{-(t-s)} d_s + c(1+r)^t \quad (19)$$

- The stock price is equal to the present discounted value of dividends (**fundamental value**) plus a **bubble component** growing at rate r

Economic example: unstable difference equation II

- Again, without further information, we cannot uniquely pin down c and therefore p_t
- Let's impose a boundary condition of non-explosive/bounded stock prices, i.e. $c = 0$:

$$p_t = \sum_{s=t+1}^{\infty} \left(\frac{1}{1+r} \right)^{-(t-s)} d_s \quad (20)$$

- In this case, the (fundamental) stock price is equal to the discounted future stream of dividends
- But: does this boundary condition make sense?

Summary

- For stable LDEs, we solve backward and need an initial condition for uniqueness
- For unstable LDEs, we solve forward and need a terminal condition for a unique and bounded solution
- This suggests two problems that can arise:
 1. if we have an initial condition paired with an unstable root, no bounded solution exists (called **no solution**)
 2. if we only have a terminal condition paired with a stable root, infinitely many bounded solutions exist (called **indeterminacy**)

LDEs: potential complications

In practical applications, there can be two complications

1. Models may involve difference equations of order higher than 1
2. Economic models usually involve both stable and unstable equations
→ we now know how to handle both of them separately, but what if the equations are simultaneously related?

Higher order linear difference equations

- The second-order linear difference equation

$$x_t = a_1 x_{t-1} + a_2 x_{t-2} + e_t \quad (21)$$

can be brought into **companion form** and thus handled as a system of two first-order equations

- Defining an auxiliary variable $z_t := x_{t-1}$ we can form a two-dimensional first-order system

$$\begin{bmatrix} x_t \\ z_t \end{bmatrix} = \begin{bmatrix} a_1 & a_2 \\ 1 & 0 \end{bmatrix} \begin{bmatrix} x_{t-1} \\ z_{t-1} \end{bmatrix} + \begin{bmatrix} 1 \\ 0 \end{bmatrix} e_t$$

- Generally, an n -th order linear difference equation can be analyzed as an n -dimensional system of first-order LDEs
- But how to deal with these systems?

The n -dimensional case

- Consider the n -dimensional first-order LDE vector system

$$x_t = Ax_{t-1} + e_t \quad (22)$$

with the square **transition matrix** A containing constant coefficients

- We assume A to be **non-singular**, i.e. invertible
- We will learn how to **decouple** the system into n independent equations that can be solved separately
- One way is to **diagonalize the transition matrix** via a **Jordan eigenvalue decomposition**
- If A has real and distinct eigenvalues, it can be diagonalized using the eigenvalue decomposition as described below
- If these assumptions are not satisfied, other methods like the generalized Schur decomposition (e.g. Klein 2000), are available

Eigenvalues and eigenvectors: definition

- The scalar λ is called an **eigenvalue** (or characteristic root) of A and the vector z a corresponding **eigenvector** if they satisfy

$$\lambda z = Az \quad (23)$$

or equivalently

$$(A - \lambda I)z = 0 \quad (24)$$

- A non-zero solution requires the matrix $A - \lambda I$ to be singular
→ determinant must be 0
- Hence, for given matrix A the solution(s) for λ can be derived as the roots of the **characteristic equation**

$$|A - \lambda I| = 0 \quad (25)$$

- The resulting characteristic equation has n solutions and A therefore n (potentially complex) eigenvalues: $\lambda = h \pm iv$.

Stable eigenvalues

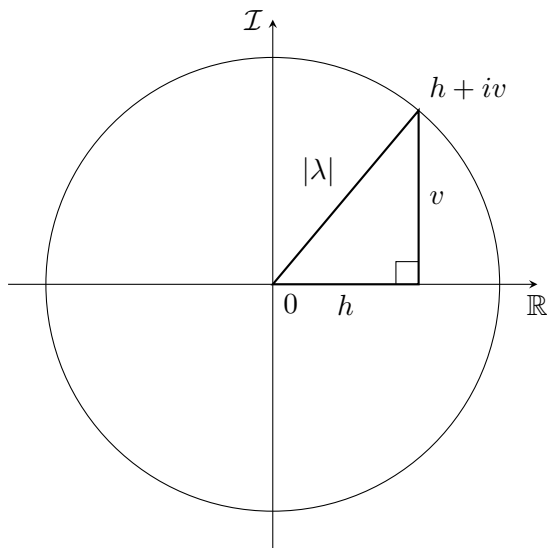
- The **modulus** of an eigenvalue is given by its Euclidean norm
$$|\lambda| = \sqrt{h^2 + v^2}$$

Definition 1

An eigenvalue is stable if its modulus is strictly smaller than one: $|\lambda| < 1$

- For now we will only consider cases with real eigenvalues
→ stable eigenvalue is one that is smaller than one in absolute value

Modulus of an eigenvalue



- In the complex plane, a modulus below 1 can be represented as being **inside of the unit circle**
- A thorny issue is the treatment of unit eigenvalues (unit roots)
- In economic problems it often makes sense to lump them together with the stable roots

Jordan Eigenvalue Decomposition

- Define Λ as the diagonal matrix with the eigenvalues of A on its main diagonal:

$$\Lambda = \begin{bmatrix} \lambda_1 & 0 & \cdots & 0 \\ 0 & \lambda_2 & \ddots & \vdots \\ \vdots & \ddots & \ddots & 0 \\ 0 & \cdots & 0 & \lambda_n \end{bmatrix} \quad (26)$$

- Let D collect the eigenvectors

$$D = [z^1, \dots, z^n] \quad (27)$$

- With distinct real eigenvalues it can be shown that A can be decomposed via the **Jordan Eigenvalue Decomposition** into

$$A = D\Lambda D^{-1} \quad (28)$$

Application to our difference equation

- Premultiplying our system (22) by D^{-1} we get

$$D^{-1}x_t = D^{-1}Ax_{t-1} + D^{-1}e_t \stackrel{(28)}{=} \Lambda D^{-1}x_{t-1} + D^{-1}e_t \quad (29)$$

Defining

$$\zeta_t = D^{-1}x_t \quad (30)$$

$$v_t = D^{-1}e_t \quad (31)$$

we can write this an uncoupled system

$$\zeta_t = \Lambda \zeta_{t-1} + v_t \quad (32)$$

Solving the decoupled system

- The ζ_{it} can be solved separately with the single equation method used above (with the initial/terminal conditions for x transformed into ones for ζ)
- The eigenvalues of A appearing on the diagonal of Λ form the coefficients in the transformed equations
 - Equations associated with an unstable eigenvalue are solved forward
 - Equations associated with a stable eigenvalue are solved backward
- Once the solution ζ_t has been found, the solution for the original variables can be computed from

$$x_t = D\zeta_t \quad (33)$$

The importance of backward-looking stochastic LDEs

- Most of the time we will be concerned with stochastic problems, giving rise to stochastic linear difference equations
- In this case, the perfect foresight assumption on e_t is replaced by e_t being random
- They can be handled very similarly as deterministic ones (as we will see later)
- Typically, solutions to economic problems take the form of backward-looking SLDEs in the **predetermined variables**, i.e. the ones that are restricted by initial conditions
- Because these predetermined variables cannot freely adjust every period, they are also called **state variables**
- In this section, we will characterize stable and unique solutions in the state variables

Stochastic LDEs: Setup

- Consider a process $\{x_t\}$ of n -dimensional vectors, recursively generated by the following first-order linear difference equation:

$$x_{t+1} = A_o x_t + C e_{t+1}, \text{ for } t = 0, 1, \dots \quad (34)$$

where

- $\{e_t\}$ for $t = 0, 1, \dots$ is a sequence of $m \times 1$ -dimensional random vectors
- x_0 is a given initial condition drawn from a distribution with mean μ_0 and covariance $\Sigma_0 = E[(x_0 - \mu_0)(x_0 - \mu_0)']$
- x_t is an $n \times 1$ **state vector** containing variables that are observed by agents at time t
- A_o is an $n \times n$ transition matrix,
- C is an $n \times m$ matrix related to the covariance
- The generated $\{x_t\}$ is then also a random sequence (**stochastic process**)
- Note the intimate connection of (34) to VARs from time series econometrics (where $n = m$)

Stochastic shock structure

- The $\{e_t\}$ process generates a sequence of information sets $\{J_t\}$ with $J_t = \{e_t, \dots, e_1, x_0\}$.
- The sequence $\{e_t\}$ is assumed to satisfy

$$E[e_{t+1}|J_t] = 0, \quad (35)$$

$$E[e_{t+1}e'_{t+1}|J_t] = I_m, \quad (36)$$

where $E_t \equiv E[\cdot|J_t]$ denotes the conditional expectations based on information set J_t

- Hence, $\{e_t\}$ is a sequence of serially uncorrelated random vectors with an unconditional mean equal to zero (shown below)
- *Exercise 5:* Show that x_t is contained in the information set J_t

Expectations operator: mathematical rules

- $E_t [X_{t+1} + Y_{t+1}] = E_t [X_{t+1}] + E_t [Y_{t+1}] .$
- In general: $E_t [X_{t+1} \cdot Y_{t+1}] \neq E_t [X_{t+1}] \cdot E_t [Y_{t+1}]$
- **Law of iterated expectations:**

$$E_t [E_{t+1} [X_{t+2}]] = E_t [X_{t+2}]$$

- Constants and variables contained in the information set:
 - $E_t [X_t] = X_t .$
 - $E_t [X_t \cdot Y_{t+1}] = X_t \cdot E_t [Y_{t+1}] .$
 - $E_t [b \cdot X_{t+1}] = b \cdot E_t [X_{t+1}] .$

- Taking derivatives

- $\frac{\partial}{\partial Y_{t+1}} E_t [f(Y_{t+1})] = E_t \left[\frac{\partial}{\partial Y_{t+1}} f(Y_{t+1}) \right] .$

- *Exercise 6:* Show that

$$E[e_t] = 0 \tag{37}$$

Conditional mean and covariance

- *Exercise 7:* Show that, if x_t follows the SLDE in (34), the optimal forecast of x_{t+1} (in the mean squared sense) for a given information set at time t is

$$E[x_{t+1}|J_t] = A_o x_t, \quad (38)$$

- The one-step ahead forecast error is

$$x_{t+1} - E_t x_{t+1} = C e_{t+1} \quad (39)$$

- The **conditional covariance matrix of x_{t+1}** (sometimes called **one-step ahead mean squared prediction error**), can be computed by postmultiplying the prediction error by its transpose and taking conditional expectations:

$$\begin{aligned} E_t \left[(x_{t+1} - E_t x_{t+1}) (x_{t+1} - E_t x_{t+1})' \right] &= E_t \left[C e_{t+1} (C e_{t+1})' \right] \\ &= E_t C e_{t+1} e_{t+1}' C' = C C'. \end{aligned} \quad (40)$$

Moving average representation

- The **autoregressive (AR)** first-order linear difference equation (34) can alternatively be written as a **moving average (MA) process**
- Exercise 8:* Show that

$$x_{t+1} = A_o x_t + C e_{t+1} = \sum_{i=1}^{t+1} A_o^{t+1-i} C e_i + A_o^{t+1} x_0 \quad (41)$$

- Leading this by $j - 1$ -periods

$$x_{t+j} = \sum_{s=t+1}^{t+j} A_o^{t+j-s} C e_s + A_o^j x_t \quad (42)$$

and applying conditional expectations yields the j -step ahead prediction

$$E_t x_{t+j} = A_o^j x_t. \quad (43)$$

- The associated j -step ahead prediction error is then given by

$$x_{t+j} - E_t x_{t+j} = \sum_{s=t+1}^{t+j} A_o^{t+j-s} C e_s = \sum_{i=0}^{j-1} A_o^i C e_{t+j-i} \quad (44)$$

j -step ahead covariance matrix

- The covariance matrix of the j -step ahead forecast error follows from

$$\begin{aligned}
 \Xi_t(j) &= E_t \left[(x_{t+j} - E_t x_{t+j}) (x_{t+j} - E_t x_{t+j})' \right] \\
 &\stackrel{(44)}{=} E_t \left[\sum_{i=0}^{j-1} A_o^i C e_{t+j-i} \left(\sum_{i=0}^{j-1} A_o^i C e_{t+j-i} \right)' \right] \\
 &= E_t \left[\sum_{i=0}^{j-1} A_o^i C e_{t+j-i} \sum_{i=0}^{j-1} e'_{t+j-i} C' (A_o^i)' \right] \\
 &= \sum_{i=0}^{j-1} A_o^i C C' (A_o^i)' ,
 \end{aligned} \tag{45}$$

where we used that $E_t(e_{t+i}e'_{t+i}) = I$ and $E_t(e_{t+i}e'_{t+k}) = 0$ for $i \neq k$ and $i, k > 0$.

- Thus, uncertainty increases in the forecast horizon j

Generic formulation with a constant

- We will consider the concept of **covariance stationarity** using a system where we have separated out constant terms:

$$A_o = \begin{bmatrix} A & \tilde{A} \\ 0 & 1 \end{bmatrix}, C = \begin{bmatrix} C_1 \\ 0 \end{bmatrix}, \quad (46)$$

where

- A is an $(n-1) \times (n-1)$ matrix with only stable eigenvalues
- \tilde{A} is an $(n-1) \times 1$ column vector
- Correspondingly partitioning $x'_t = [x'_{1t} \quad x'_{2t}]$, equation (34) can be written as

$$x_{1t+1} = Ax_{1t} + \tilde{A}x_{2t} + C_1e_{t+1}, \quad (47)$$

$$x_{2t+1} = x_{2t} \quad (48)$$

- Given that x_{2t} is constant over time (it exhibits a unit root), and thus equal to x_{20} , we define

$$B \equiv \tilde{A}x_{20} \quad (49)$$

Covariance stationarity

Definition 2

A stochastic process $\{x_t\}$ is covariance stationary if the unconditional mean is time invariant

$$Ex_t = Ex_0 \quad \forall t \quad (50)$$

and the sequence of unconditional autocovariance matrices

$$\Sigma(j) \equiv E(x_{t+j} - Ex_{t+j})(x_t - Ex_t)' \quad (51)$$

is finite and only depends on the separation between dates j , but not on t .

- Note that conditional heteroscedasticity is allowed, because it refers to conditional moments
- Our sequence $\{x_t\}$ will be covariance stationary, if all eigenvalues of A (i.e. not associated with the constant) have moduli less than one

The goal: a fixed point

- We will be looking for an initial distribution with mean $\mu_0 = E x_0$ and covariance $\Sigma_0 = E [(x_0 - \mu_0)(x_0 - \mu_0)']$ that, if x_0 is drawn from this distribution, makes x_t covariance stationary
- If instead starting from an arbitrary initial value x_0 , the process will still converge to this stationary distribution as $t \rightarrow \infty$

The unconditional mean

- Taking **unconditional expectations** on both sides of (47), yields

$$Ex_{1t+1} = AEx_{1t} + B. \quad (52)$$

- If μ_0 is the **stationary value** of Ex_{1t} , it needs to satisfy

$$\mu_0 = A\mu_0 + B \Leftrightarrow \mu_0 = (I - A)^{-1}B \quad (53)$$

- As before, using that the eigenvalues are stable, we know that

$$(I - A)^{-1} = I + A + A^2 + A^3 + \dots \quad (54)$$

Convergence to the unconditional mean

- From the eigenvalue decomposition $A = D\Lambda D^{-1}$, the j^{th} power of A can be written as

$$A^j = (D\Lambda D^{-1}) (D\Lambda D^{-1}) \dots (D\Lambda D^{-1}) = D\Lambda^j D^{-1} \quad (55)$$

- Thus, the stationary mean μ_0 satisfies

$$\mu_0 = (I - A)^{-1}B \stackrel{(54)}{=} \sum_{i=0}^{\infty} A^i B \stackrel{(55)}{=} D \left(\sum_{i=0}^{\infty} \Lambda^i \right) D^{-1} B \quad (56)$$

- If all eigenvalues of A are stable, μ_0 will have a finite value
- Exercise 9:** Show that in this case, x_{1t} will converge to the stationary value μ_0 for any initial value of x_{10}

The unconditional covariance

- Subtracting (52) from (47) yields the uncond. period t forecast error

$$(x_{1t} - Ex_{1t}) = A(x_{1t-1} - Ex_{1t-1}) + C_1 e_t \quad (57)$$

- The uncond. forecast error covariance at t , Σ_t , then satisfies

$$\begin{aligned} \Sigma_t &= E \left[(x_{1t} - Ex_{1t}) (x_{1t} - Ex_{1t})' \right] \\ &\stackrel{(57)}{=} E \left[(A(x_{1t-1} - Ex_{1t-1}) + C_1 e_t) (A(x_{1t-1} - Ex_{1t-1}) + C_1 e_t)' \right] \\ &= E \left[A(x_{1t-1} - Ex_{1t-1}) (x_{1t-1} - Ex_{1t-1})' A' + C_1 e_t e_t' C_1' \right. \\ &\quad \left. + C_1 e_t (x_{1t-1} - Ex_{1t-1})' A' + A(x_{1t-1} - Ex_{1t-1}) e_t' C_1' \right] \\ &= AE \left[(x_{1t-1} - Ex_{1t-1}) (x_{1t-1} - Ex_{1t-1})' \right] A' + E \left[C_1 e_t e_t' C_1' \right] \\ &= A \Sigma_{t-1} A' + C_1 C_1', \end{aligned} \quad (58)$$

where the second-to-last line uses that e_t is orthogonal to $x_{1t-1} - Ex_{1t-1}$

The unconditional covariance II

- The result is a **discrete Lyapunov** equation

$$\Sigma_t = A\Sigma_{t-1}A' + C_1C_1' \quad (58)$$

- A stationary value Σ_0 for the covariance matrix of x_{1t} has to satisfy (with $Ex_{1t} = \mu_0$)

$$\Sigma_0 = A\Sigma_0A' + C_1C_1' \quad (59)$$

- Iterating backwards yields the solution

$$\Sigma_0 = \sum_{k=0}^{\infty} A^k C_1 C_1' A^{k'} \quad (60)$$

- Exercise 10:* Verify that this solves (59) if A is a stable matrix
- Comparing the j -step ahead forecast error variance, (45), with the unconditional one in (60) reveals that the unconditional covariance is the limit of the former

Inspecting the stationary covariance matrix

- If the eigenvalues of A have moduli less than one, the unconditional covariance matrix Σ_0 takes a finite value:

$$\begin{aligned}
 \Sigma_0 &\stackrel{(55)}{=} \sum_{j=0}^{\infty} D \Lambda^j D^{-1} C_1 C_1' \left(D \Lambda^j D^{-1} \right)' \\
 &= D \sum_{j=0}^{\infty} \Lambda^j \left[D^{-1} C_1 \right] \left[C_1' (D^{-1})' \right] \left(\Lambda^j \right)' D' \\
 &= D \left(\sum_{j=0}^{\infty} \Lambda^j \tilde{C} \tilde{C}' \left(\Lambda^j \right)' \right) D',
 \end{aligned} \tag{61}$$

where $\tilde{C} = D^{-1} C_1$ and $\tilde{C}' = C_1' (D^{-1})'$

- Thus, if the covariance matrix of x_{10} is Σ_0 , it will remain constant over time
- Otherwise, the sequence of covariance matrices will asymptotically converge to Σ_0

The unconditional autocovariances

- This shows that the contemporaneous covariance at lag 0, $\Sigma_0 = \Sigma_0(0)$, only depends on j , but not on t
- But what about other leads and lags?
- The unconditional forecast error at time $t + j$ satisfies

$$(x_{1t+j} - Ex_{1t+j}) \stackrel{(44)}{=} A^{j+1} (x_{1t-1} - Ex_{1t-1}) + C_1 e_{t+j} + \dots + A^j C_1 e_t \quad (62)$$

- *Exercise 11*: Show that the autocovariance at lead/lag j

$$\Sigma_t(j) \equiv E(x_{1t+j} - Ex_{1t+j})(x_{1t} - Ex_{1t})' \quad (63)$$

is given by

$$\Sigma_t(j) = A^j \Sigma_t(0) \quad (64)$$

- Thus, if $\Sigma_t(0)$ does not depend on t , neither does $\Sigma_t(j)$

State space systems

- We have already characterized the **state transition equation**, i.e. the evolution of the predetermined state variables x_t over time

$$x_{t+1} = A_o x_t + C e_{t+1}, \text{ for } t = 0, 1, \dots \quad (34)$$

- But often we are also interested in functions of the state variables:

$$y_t = G x_t, \quad (65)$$

where y_t is a $n_y \times 1$ vector and G a $n_y \times n_x$ matrix

- Equation (65) linking endogenous non-state variables to state variables is called a **measurement equation/observation equation**
- Equations (34) and (65) together form a **state-space system**, describing the evolution of all variables of interest

Moving average representations (again)

- We already derived the moving average representation for the state variables

$$x_{t+1} = \sum_{i=1}^{t+1} A_o^{t+1-i} C e_i + A_o^{t+1} x_0 \quad (41)$$

- From (65) then follows

$$y_t = G \sum_{i=0}^t A_o^{t-i} C e_i + G A_o^t x_0 \quad (66)$$

- *Exercise 12*: Show that

$$E[y_t | x_0] = G A_o^t x_0 \quad (67)$$

and using (45) that

$$E_t \left[(y_{t+j} - E_t y_{t+j}) (y_{t+j} - E_t y_{t+j})' \right] = G \left[\sum_{i=0}^{j-1} A_o^i C C' (A_o^i)' \right] G' \quad (68)$$

Impulse response functions

- We will often be concerned with the dynamic response of a variable y_{t+h} to a shock ε_t occurring at time t :

$$\frac{\partial y_{t+h}}{\partial \varepsilon_t} \text{ for some } t \text{ and some } h \text{ (typically } \geq 0) \quad (69)$$

- The moving average representations (41) and (66) reveal that the **impulse-response functions** for $h \geq 0$ are given by

$$\frac{\partial y_{t+h}}{\partial \varepsilon_t} = GA_o^h C \quad (70)$$

$$\frac{\partial x_{t+h}}{\partial \varepsilon_t} = A_o^h C \quad (71)$$

- If all eigenvalues are inside of the unit circle, the effect of shocks will die out over time

Outline

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 - 4.1 Lagrangian formulations
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Introduction

- We have made great headway in solving and understanding LDEs, but we still do not know where the LDEs come from in economic applications
- They follow from economic optimization problems in discrete time
- We again start with the deterministic case, before turning to the stochastic case

Setup: the sequence problem

1. Choose an infinite sequence of **control variables**, $\{u_t\}_{t=0}^{\infty}$
2. to maximize a discounted time-separable objective

$$\max_{\{u\}_{t=0}^{\infty}} \sum_{t=0}^{\infty} \beta^t r(x_t, u_t), \quad (72)$$

where $\beta \in (0, 1)$ is a discount factor and r the instantaneous return function,

3. subject to a set of equations constraining the evolution of the **state variables** $\{x_t\}_{t=0}^{\infty}$

$$x_{t+1} \leq g_t(x_t, u_t) \quad (73)$$

4. for a given initial value x_0

Classification of variables

- When discussing LDEs, we already saw the important distinction between predetermined variables and non-predetermined ones
- In economic problems, we classify variables into endogenous and exogenous state variables and control variables
- **State variables:** cannot be changed by the agent in the current period, because they are **predetermined**. They characterize the state of the system at any point in time.
 - **exogenous** state: cannot be changed by the agent at any time, but follow exogenous (stochastic) law of motion
 - **endogenous** state: are predetermined at time t , but can be changed by the agent at time $t + j, j > 0$
- **Control variables:** all variables that the agent can change at time t (can be an endogenous state variable at $t + j, j > 0$)
- We will discuss examples later on

The goal

- Derive a (time-invariant) **policy function** h that maps the state x_t into the control u_t such that the sequence of controls generated by the two functions

$$u_t = h(x_t) \tag{74}$$

$$x_{t+1} \leq g(x_t, u_t), \tag{73}$$

starting from the initial condition x_0 at $t = 0$, solves the problem

- Knowing $h : D \rightarrow \mathbb{R}$ is equivalent to knowing $\{u_t\}_{t=0}^{\infty}$
- Note that both $\{u_t\}_{t=0}^{\infty}$ and h are infinite-dimensional objects
- The structure of the system (73)-(74) is called **recursive**: starting with x_0 , we can generate the sequence for $\{u_t, x_{t+1}\}$ by iteratively plugging in
- We will only encounter functions, but h could be a **correspondence**

Getting the solution

- There are two different solution techniques:
 1. Dynamic programming (covered in Advanced Macro II)
 2. Lagrange-based methods
- Lagrange is typically more intuitive, but requires differentiability
- Dynamic programming works with functions instead of sequences, requiring bigger formal investments, but
 - it has computational advantages, particularly when the problem is not differentiable
 - solution characteristics like existence, uniqueness, boundedness, etc. are often easier to derive

The Lagrangian formulation

- If everything is nicely differentiable, discrete time optimization problems can also be solved by applying the **Lagrange formulation** instead of dynamic programming via Bellman equations
- The Lagrangian sequence problem formulation of the original problem is

$$\mathcal{L} = \sum_{t=0}^{\infty} \beta^t [r(x_t, u_t) + \lambda'_t (g(x_t, u_t) - x_{t+1})], \quad (75)$$

where x is an n -dimensional vector of states, λ_t is an n -dimensional vector of Lagrange multipliers, and x_0 is given

- By convention: discounting applies to the Lagrange multiplier as well
- We seek to maximize the Lagrange function by choosing sequences $\{u_t, x_{t+1}\}_{t=0}^{\infty}$ (and λ_t)
- Note that the Lagrange multiplier part in (75) is equivalent to $-\lambda'_t (x_{t+1} - g(x_t, u_t))$

The first order conditions

- The first-order necessary conditions at every point in time t are given by the transversality constraint (TVC) and

$$0 = \beta^t \frac{\partial r(x_t, u_t)}{\partial u_t} + \beta^t \lambda_t \frac{\partial g(x_t, u_t)}{\partial u_t} \quad (76)$$

$$0 = -\beta^t \lambda_t + \beta^{t+1} \frac{\partial r(x_{t+1}, u_{t+1})}{\partial x_{t+1}} + \beta^{t+1} \lambda_{t+1} \frac{\partial g(x_{t+1}, u_{t+1})}{\partial x_{t+1}} \quad (77)$$

$$0 = g(x_t, u_t) - x_{t+1}, \quad (78)$$

where we used the vector calculus rule that $\frac{\partial \lambda' a}{\partial a} = \lambda$ (see e.g. Greene 2011, p.1009)

- Due to the stationary problem, the FOCs are time-invariant
- Simplifying gives

$$\frac{\partial r(x_t, u_t)}{\partial u_t} + \lambda_t \frac{\partial g(x_t, u_t)}{\partial u_t} = 0 \quad (79)$$

$$\beta \frac{\partial r(x_{t+1}, u_{t+1})}{\partial x_{t+1}} + \beta \lambda_{t+1} \frac{\partial g(x_{t+1}, u_{t+1})}{\partial x_{t+1}} = \lambda_t \quad (80)$$

The transversality condition

- Starting from a finite horizon problem, if we let $T \rightarrow \infty$, the terminal condition for capital leads to the **transversality condition**

$$\lim_{T \rightarrow \infty} \beta^T \lambda_{1,T} k_{T+1} = 0 \quad (81)$$

- The transversality condition (81) is a necessary condition for the maximization problem
- We will repeatedly face infinite horizon problems leading to this type of condition
- This TVC essentially implies that the discounted value of wealth (i.e., capital) in terms of utility has to be zero for the limiting case where time approaches infinity
- TVCs serve as terminal conditions that have to be taken into account when solving the set of first-order conditions (which will typically take the form of vector difference equations)

The neoclassical growth model

- Consider a Ramsey (1928)-Cass (1965)-Koopmans (1965)-type model with one representative household, supplying labor inelastically:

$$N = 1$$

- Felicity is logarithmic in consumption:

$$u(C_t) = \ln C_t \quad (82)$$

and households have discount factor $\beta \in [0, 1)$

- Production technology is Cobb-Douglas with capital share $\alpha \in (0, 1)$:

$$Y_t = F(K_t, N_t) = AK_t^\alpha N_t^{1-\alpha}, \quad (83)$$

with TFP $A > 0$ being exogenous

- The law of motion for capital is given by

$$K_{t+1} = (1 - \delta)K_t + I_t, \quad (84)$$

where $\delta \in (0, 1]$ is the depreciation rate

- We assume a closed economy, i.e. the resource constraint is

$$Y_t = C_t + I_t \quad (85)$$

Stock notations

- The capital stock is a predetermined variable
- There are two timing conventions in the literature:
 1. the **stock at the end of period notation** uses the capital stock at the end of the period when investment has already been added:

$$K_t = (1 - \delta)K_{t-1} + I_t \quad (86)$$

$$Y_t = AK_{t-1}^\alpha N_t^{1-\alpha} \quad (87)$$

2. the **stock at the beginning of period notation** uses the capital stock at the beginning of the current period, before current investment has been added:

$$K_{t+1} = (1 - \delta)K_t + I_t \quad (88)$$

$$Y_t = AK_t^\alpha N_t^{1-\alpha} \quad (89)$$

- The **stock at the end of period notation** is typically more intuitive, because it endows the capital stock with the timing at which it is decided and makes the predeterminedness explicit

The planner's problem

- Define lower-case letters as per capita variables, i.e. $x_t = X_t/N_t$
- *Exercise 13*: Assuming full depreciation $\delta = 1$, show that

$$k_{t+1} = Ak_t^\alpha - c_t \quad (90)$$

- The social planner's program then is:

$$\begin{aligned} \max_{\{c_t\}_{t=0}^{\infty}} \sum_{t=0}^{\infty} \beta^t \ln c_t \\ \text{s.t. } k_{t+1} = Ak_t^\alpha - c_t, \\ k_0 > 0 \text{ and TVC given} \end{aligned} \quad (90)$$

- Variable classification
 - State variables: k_t
 - Control variables: c_t, k_{t+1}

The Brock and Mirman (1972)-model

- For the simple neoclassical growth model from above, the Lagrange sequence problem reads

$$\mathcal{L} = \sum_{t=0}^{\infty} \beta^t [\ln c_t + \lambda_t (Ak_t^\alpha - c_t - k_{t+1})]. \quad (91)$$

- The first-order conditions (79)-(80) are given by

$$\lambda_t = c_t^{-1}, \quad (92)$$

$$\lambda_t = \beta \lambda_{t+1} \alpha A k_{t+1}^{\alpha-1}, \quad (93)$$

where we used that

$$\frac{\partial r(x_{t+1}, u_{t+1})}{\partial x_{t+1}} = \frac{\partial \ln c_{t+1}}{\partial k_{t+1}} = 0 \quad (94)$$

- Eliminating the Lagrange multiplier yields the Euler equation:

$$\frac{c_{t+1}}{c_t} = \beta \alpha A k_{t+1}^{\alpha-1} \quad (95)$$

Reminder: The Keynes-Ramsey rule I

- The Euler, together with the transition law (90), implicitly characterizes the optimal sequence of c_t and k_t for a given initial value k_0 (second order difference equation)
- The Euler equation for consumption per worker yields the **Keynes-Ramsey rule**

$$\frac{\Delta c_{t+1}}{c_t} \approx \log\left(\frac{c_{t+1}}{c_t}\right) = \frac{r_{t+1,t} - \rho}{\sigma} = (r_{t+1,t} - \rho) \times EOIS \quad (96)$$

where we used:

- the approximation $\log(1+x) \approx x$ for small x
- $\beta = 1/(1+\rho)$, where ρ is the rate of time preference
- CRRA preferences of the type $(C_t^{1-\sigma} - 1)/(1-\sigma)$ with risk aversion coefficient σ so that the **intertemporal elasticity of substitution** is $EOIS = 1/\sigma$
- $r_{t+1,t} = f'(k_{t+1}) + (1-\delta) = \alpha A k_{t+1}^{\alpha-1}$

Reminder: The Keynes-Ramsey rule II

- The Euler equation for consumption per worker yields the **Keynes-Ramsey rule**

$$\frac{\Delta c_{t+1}}{c_t} \approx \log \left(\frac{c_{t+1}}{c_t} \right) = \frac{r_{t+1,t} - \rho}{\sigma} = (r_{t+1,t} - \rho) \times EOIS \quad (96)$$

- If $r > \rho$: Giving up one unit of consumption today and saving it yields r , but the HH only discounts with $\rho \Rightarrow$ HH shifts consumption to the future, i.e. consumption path is increasing
- Marginal utility in the future must be lower and consumption higher than today to make RHS of (95) equal to LHS, given $r_{t+1,t} - \rho > 0$
- In EQ: shifting consumption in any direction at the margin does not increase discounted marginal utility
- The willingness to shift consumption intertemporally depends on the EOIS $1/\sigma$: the higher the EOIS the more consumption responds to a given $r - \rho$

The method of undetermined coefficients

- We are left with a problem: the Euler equation only implicitly characterizes the solution
- In general, it is impossible to immediately obtain the policy functions from these equilibrium conditions
- However: the Brock and Mirman (1972)-model due to its log-linear structure has a simple analytical solution that can be guessed and verified via the **method of undetermined coefficients**
- Idea: guess a functional form for h but leave the coefficients unknown; functional form should only depend on the state of the system and be similar to the model structure
- Plug this guess (repeatedly) into the FOCs and try to solve for the unknown coefficient(s)
- If we can find a value for the coefficients (i.e. a function of only parameters, not variables) that satisfies the FOCs, we have found one (of potentially many) solution

Getting the solution

- *Exercise 14*: Use the guess of a constant savings rate

$$k_{t+1} = \phi A k_t^\alpha \tag{97}$$

and show that $\phi = \alpha\beta$ solves the problem

Intro

- This section presents methods of dynamic optimization under **uncertainty**, which will repeatedly be applied in this course

Stochastic structure

- We introduce random events $s_t \in S$ where S is the finite set of possible events $S = \{\sigma_a, \sigma_b, \sigma_c, \dots, \sigma_n\}$ that can happen in any period t
- The unconditional probability of an event is given by $\pi(s_t)$ and

$$\sum_{s_t \in S} \pi(s_t) = 1 \quad (98)$$

- The **event history** is given by $s^t = (s_t, s_{t-1}, \dots, s_0)$ and summarizes the realizations of all events from period 0 to period t .
- The set of event histories is then $S^t = S \times S \times \dots \times S$, with $s^t \in S^t$
- The probability of a particular event history s^t is given by $\pi(s^t)$, where $\pi(s^t) > 0$ for all $s^t \in S^t$ and t
- We could relax the assumption that the same set of events S can realize with some strictly positive probability in each period, but
 - notation would be heavier
 - problem would not be time-invariant

Probability of event history

- The probability of a particular event history $\pi(s^t) = \pi(s_t, s_{t-1}, \dots, s_0)$ is the joint probability of the events s_t, \dots, s_0 , where $\pi(s^0) = 1$ for a given initial state s^0
- Denote the conditional probability of state s_t and thus s^t given the history s^{t-1} at date $t - 1$ with $\pi(s^t | s^{t-1})$
- The joint probability can be factored as

$$\begin{aligned}\pi(s^t) &= \pi(s^t | s^{t-1}) \cdot \pi(s^{t-1} | s^{t-2}) \cdot \dots \cdot \pi(s^1 | s^0) \cdot \pi(s^0) \\ &= \pi(s_t | s^{t-1}) \cdot \pi(s_{t-1} | s^{t-2}) \cdot \dots \cdot \pi(s_1 | s^0),\end{aligned}\tag{99}$$

where we used that $\pi(s^0) = 1$.

Markov processes

- In general the probability of a particular event s_t occurring at time t depends on the complete history of past events
- That makes general processes hard to work with, because past events become states that need to be tracked
- Life becomes easier when working with Markov processes that only require keeping track of some history

Definition 3 (n^{th} -order Markov process)

s_t is generated by an n^{th} -**order Markov process** if the distribution of s_t conditional on n lags is the same as the distribution of s_t conditional on all lags:

$$\pi(s_t | s^{t-1}) = \pi(s_t | s_{t-1}, \dots, s_{t-n}) \quad (100)$$

First-order Markov processes

- A first-order Markov process is then characterized by

$$\pi(s_t | s_{t-1}) = \pi(s_t | s^{t-1}) \quad (101)$$

- The probability of a particular event only depends on the last event, but not on the entire history of earlier events
- The probability of event history s^t ,

$$\pi(s^t) = \pi(s_t, s_{t-1}, \dots, s_0) \quad (102)$$

for a given initial state s_0 then simplifies to

$$\begin{aligned} \pi(s^t) &= \pi(s_t | s^{t-1}) \cdot \pi(s_{t-1} | s^{t-2}) \cdot \dots \cdot \pi(s_1 | s^0) \\ &= \pi(s_t | s_{t-1}) \cdot \pi(s_{t-1} | s_{t-2}) \cdot \dots \cdot \pi(s_1 | s_0) \end{aligned} \quad (103)$$

Example

- Consider a 2-state Markov process: each period the event is $s_t \in S = \{\sigma_a, \sigma_b\}$ with $s_0 = \sigma$ being a given fixed event $\in S$
- The history of events up to t is then

$$s^t = (s_t, s_{t-1}, \dots, s_0) \in S^t = S_t \times S_{t-1} \times \dots \times S_0 \quad (104)$$

- Thus, for the periods $t = 0, 1, 2$ the sets of event histories are

$$S_0 = \sigma \quad (105)$$

$$S^1 = \{(\sigma_a, \sigma), (\sigma_b, \sigma)\} \quad (106)$$

$$S^2 = \{(\sigma_a, \sigma_a, \sigma), (\sigma_a, \sigma_b, \sigma), (\sigma_b, \sigma_a, \sigma), (\sigma_b, \sigma_b, \sigma)\} \quad (107)$$

- The unconditional probability for event history $\pi(s^2)$ is given by

$$\begin{aligned} \pi(s^2) &= \pi(s_2|s^1) \cdot \pi(s_1|s^0) \cdot \pi(s^0) \\ &= \pi(s_2|s_1) \cdot \pi(s_1|s_0) \cdot \pi(s_0) \end{aligned} \quad (108)$$

Example continued

- The unconditional probability for the particular history $s^2 = (\sigma_b, \sigma_a, \sigma)$ is given by

$$\pi((\sigma_b, \sigma_a, \sigma)) = \pi(\sigma_b | \sigma_a) \cdot \pi(\sigma_a | \sigma), \quad (109)$$

where we used that $\pi(s^0) = \pi(\sigma) = 1$.

A stochastic concave programming problem

- Consider a problem with a time-separable differentiable and concave return function $r(x_t, y_t)$ and discount factor $\beta \in (0, 1)$
- The sequence of control variables y_t is chosen to maximize an intertemporal objective, subject to a set of constraints on the evolution of the endogenous state variables x_t

$$x_{t+1} = g(x_t, y_t) \quad (110)$$

- The set $\{(x_{t+1}, x_t) : x_{t+1} \leq g(x_t, y_t)\}$ is convex and compact for admissible y_t
- We assume events s_t realize at the beginning of each period, i.e. after y_{t-1} has been chosen, but before y_t
- Controls y_t will generally depend on event history s^t , i.e. $y_t(s^t)$, while state variables x_t will depend on history s^{t-1} , i.e. $x_t(s^{t-1})$
- Note that they might depend on more than the random events

Stochastic maximization problem

- Stochastic problems typically involve the maximization of an **expected** sum of discounted returns (e.g. **von Neumann-Morgenstern utility**), subject to set of constraints that has to hold for each state of the world:

$$\max_{\{y_t(s^t), x_{t+1}(s^t)\}_{t=0}^{\infty}} \sum_{t=0}^{\infty} \sum_{s^t \in S^t} \beta^t \pi(s^t) r(x_t(s^{t-1}), y_t(s^t)), \quad (111)$$

$$\text{s.t.} \quad x_{t+1}(s^t) = g(x_t(s^{t-1}), y_t(s^t), s_t),$$

where initial values x_0 and s_0 are given and random events s are generated by a first-order Markov process

- The Lagrangian for this problem is given by

$$\begin{aligned} \mathcal{L}(x_0, s_0) = & \sum_{t=0}^{\infty} \sum_{s^t \in S^t} \beta^t \pi(s^t) \left[r(x_t(s^{t-1}), y_t(s^t)) \right. \\ & \left. + \lambda'_t(s^t) \left(g(x_t(s^{t-1}), y_t(s^t), s_t) - x_{t+1}(s^t) \right) \right] \end{aligned} \quad (112)$$

The first order conditions

- The first-order conditions with respect to $y_t(s^t) \forall s^t, s^{t+1} \in S^t$ are

$$0 = \beta^t \pi(s^t) \frac{\partial r(x_t(s^{t-1}), y_t(s^t))}{\partial y_t(s^t)} + \beta^t \pi(s^t) \lambda_t(s^t) \frac{\partial g(x_t(s^{t-1}), y_t(s^t), s_t)}{\partial y_t(s^t)}, \quad (113)$$

- The FOCs with respect to $x_{t+1}(s^t)$ are

$$\begin{aligned} \pi(s^t) \lambda_t(s^t) = & \sum_{s^{t+1}|s^t} \beta \pi(s^{t+1}) \frac{\partial r((x_{t+1}(s^t), y_{t+1}(s^{t+1})))}{\partial x_{t+1}(s^t)} \\ & + \sum_{s^{t+1}|s^t} \beta \pi(s^{t+1}) \lambda_{t+1}(s^{t+1}) \frac{\partial g(x_{t+1}(s^t), y_{t+1}(s^{t+1}), s_{t+1})}{\partial x_{t+1}(s^t)}, \end{aligned} \quad (114)$$

where we sum over all possible histories s^{t+1} , conditional on the realization of a particular history s^t

- In addition, the FOC w.r.t. to λ , i.e. the constraint, has to hold

History-dependence vs. time-invariance

- Solutions for y_t and x_{t+1} that satisfy the two sets of conditions, (113)-(114) and the set of constraints

$$x_{t+1}(s^t) = g\left(x_t(s^{t-1}), y_t(s^t), s_t\right) \quad (115)$$

are generally **history-dependent**

- That is, objects are in principle time-varying functions of the history of events s^t
- However, we aim at deriving time-invariant functions for y_t and x_{t+1} that do not depend on the entire history, but on a limited set of relevant information.
- At each point in time this information is summarized in the **state**, which includes the random event s_t , i.e. the **exogenous state** and the **endogenous state** x_t

Recursive solutions

- Suppose there exists a solution for the current control y_t that can be written as a time-invariant function of the current endogenous state x_t and the current event s_t rather than of the entire history s^t .

$$y_t = h(x_t, s_t) \quad (116)$$

- In this case, the constraint requires the endogenous state to satisfy

$$x_{t+1} = g(x_t, y_t, s_t) = g(x_t, h(x_t, s_t), s_t) \quad (117)$$

- The endogenous state, like the exogenous state, is then generated by a first-order Markov process
→ the solutions for y_t and x_{t+1} inherit the Markov property

Rewriting the FOCs

- We can use the Markov property to rewrite the first order conditions (113) as

$$0 = \frac{\partial r(x_t, y_t(x_t, s_t))}{\partial y_t(x_t, s_t)} + \lambda_t(x_t, s_t) \frac{\partial g(x_t, y_t(x_t, s_t), s_t)}{\partial y_t(x_t, s_t)} \quad (113')$$

→ λ_t is also a function of x_t and s_t .

- Using the Markov property and $\frac{\pi(s^{t+1})}{\pi(s^t)} = \pi(s_{t+1}|s_t)$, we can write (114) as

$$\begin{aligned} \lambda_t(s_t) = & \sum_{s_{t+1}|s_t} \beta \pi(s_{t+1}|s_t) \frac{\partial r(x_{t+1}, y_{t+1}(x_{t+1}, s_{t+1}))}{\partial x_{t+1}} \\ & + \sum_{s_{t+1}|s_t} \beta \pi(s_{t+1}|s_t) \lambda_{t+1}(x_{t+1}, s_{t+1}) \frac{\partial g(x_{t+1}, y_{t+1}(x_{t+1}, s_{t+1}), s_{t+1})}{\partial x_{t+1}} \end{aligned} \quad (114')$$

Rewriting the FOCs II

- Introducing the expectations operator E_t conditional on information in period t (for any vector z of random variables)

$$E_t z_{t+1} = \sum_{s_{t+1}|s_t} \pi(s_{t+1}|s_t) z_{t+1}(s_{t+1}), \quad (118)$$

allows to rewrite the conditions as

$$0 = \frac{\partial r(x_t, y_t)}{\partial y_t} + \lambda_t \frac{\partial g(x_t, y_t, s_t)}{\partial y_t} \quad (113'')$$

$$\lambda_t = \beta E_t \frac{\partial r(x_{t+1}, y_{t+1})}{\partial x_{t+1}} + \beta E_t \lambda_{t+1} \frac{\partial g(x_{t+1}, y_{t+1}, s_{t+1})}{\partial x_{t+1}} \quad (114'')$$

$$x_{t+1} = g(x_t, y_t, s_t) \quad (117)$$

Inspecting the solution

- These conditions have to be satisfied by sequences $\{y_t\}_{t=0}^{\infty}$ and $\{x_{t+1}\}_{t=0}^{\infty}$, for a given sequence for s_t and x_0
- If such solutions exist, y_t and x_{t+1} are time-invariant functions of the current states x_t and s_t , i.e. identical functions $\forall t \geq 0$
- Note that y_t and x_{t+1} might further depend on the properties of the **distribution** of the random event
- If the system is linear, only the first moment of s_t will affect the solution (**certainty equivalence**)
- Otherwise, higher moments might also affect the solution

A stochastic Brock and Mirman (1972) model

- Consider the planner problem for a Brock and Mirman (1972) model with a random exogenous state variable s_t generated by a first order finite state Markov process

$$\begin{aligned} & \max_{\{c_t(s^t), k_{t+1}(s^t)\}_{t=0}^{\infty}} \sum_{t=0}^{\infty} \sum_{s^t \in S^t} \beta^t \pi(s^t) u(c_t(s^t)) \\ \text{s.t.} \quad & k_{t+1}(s^t) + c_t(s^t) = f(k_t(s^{t-1}), s_t), \end{aligned} \tag{119}$$

where u denotes a standard concave utility function, c consumption of a representative household, and k capital per capita

- The production function f exhibits neoclassical properties and output (per capita) depends on a random event via stochastic total factor productivity

The Lagrangian sequence problem

- The Lagrangian sequence problem is then given by

$$\begin{aligned} \mathcal{L} = & \sum_{t=0}^{\infty} \sum_{s^t \in S^t} \beta^t \pi(s^t) u(c_t(s^t)) \\ & + \sum_{t=0}^{\infty} \sum_{s^t \in S^t} \beta^t \pi(s^t) \lambda_t(s^t) \left(f(k_t(s^{t-1}), s_t) - k_{t+1}(s^t) - c_t(s^t) \right) \end{aligned} \quad (120)$$

- The FOCs for c_t and k_{t+1} for given histories s^t and s^{t+1} are

$$\pi(s^t) u_c(c_t(s^t)) - \pi(s^t) \lambda_t(s^t) = 0 \quad (121)$$

$$\pi(s^t) \lambda_t(s^t) - \sum_{s^{t+1} | s^t} \beta \pi(s^{t+1}) \lambda_{t+1}(s^{t+1}) f'(k_{t+1}(s^t), s_{t+1}) = 0 \quad (122)$$

Deriving the Euler equation

- The FOCs (121)-(122) can be simplified to

$$u_c(c_t(s^t)) = \lambda_t(s^t) \quad (123)$$

$$\lambda_t(s^t) = \sum_{s^{t+1}|s^t} \beta \frac{\pi(s^{t+1})}{\pi(s^t)} \lambda_{t+1}(s^{t+1}) f'(k_{t+1}(s^t), s_{t+1}) \quad (124)$$

- Now use that s_t follows a Markov process $\frac{\pi(s^{t+1})}{\pi(s^t)} = \pi(s_{t+1}|s_t)$ and suppose there exists a solution that can be written as functions of the current states, but does not depend on the previous history of events
- Then we end up with the Euler equation

$$u_c(c_t(s_t)) = \sum_{s_{t+1}|s_t} \beta \pi(s_{t+1}|s_t) u_c(c_{t+1}(s_{t+1})) f'(g(k_t, s_t), s_{t+1}) \quad (125)$$

or more compactly

$$u_c(c_t(s_t)) = E_t \beta [u_c(c_{t+1}(s_{t+1})) f'(g(k_t, s_t), s_{t+1})] \quad (126)$$

The equilibrium

- The Euler equation implies that $c_t(s_t)$ depends on current states (s_t, k_t)
- Consumption will also in general depend on the first and higher moments of the random variable s_t
- Given that consumption does not depend on the history of states, we can write the constraint as

$$k_{t+1}(k_t, s_t) = f(k_t, s_t) - c_t(s_t, k_t), \quad (127)$$

showing that capital is a function of the current states k_t and s_t

- We have demonstrated that a recursive solution might exist
- But we have not shown that it actually exists or whether it is unique
- In principle, there might even be non-recursive solutions to the problem, leading to a non-Markov evolution of the state
- We will disregard these types of solutions throughout the course

Certainty Equivalence I

- *Exercise 15*: Consider the stochastic Brock and Mirman (1972) model with log utility and stochastic TFP A_t :

$$y_t = A_t k_t^\alpha \quad (128)$$

Show that the solution is given by

$$k_{t+1} = \alpha\beta A_t k_t^\alpha \quad (129)$$

- Due to log utility, full depreciation, and Cobb-Douglas: distribution of shock process does not affect solution
- Regarding the variance, a mean-preserving spread to A_t would not alter the agents' choice
- **Certainty equivalence**: people make decisions under uncertainty by acting as if future stochastic variables were sure to turn out equal to their expected values (Obstfeld and Rogoff 1996, p. 81)
- This also implies the absence of **precautionary behavior**, because people act as if the expected value would occur with 100% certainty

Certainty Equivalence II

- As in this special case, certainty equivalence is a **property of the model**
- But: certainty equivalence can also be an **artifact of a particular solution technique**, although the underlying model does not feature it
- For example, Jensen's inequality effects cannot arise with linear solutions (more on this later)

Equilibria with several agents

- In the Brock and Mirman (1972)-model we had a central planner directly choosing the state
- But: in most macroeconomic models, the law of motion for the relevant endogenous state variables will not only depend on the decisions of one agent
- Agents will e.g. interact via markets where they are price-takers
- We will thus require an equilibrium concept where the plans of different agents are combined in a consistent way

Setup

- Let x_t be the vector of state variables under the control of the agent and X_t the vector of the same variables resulting from “markets”
- Denote with Z_t the vector of exogenous states
- The agents problem then takes the form

$$V(x_t, X_t, Z_t) = \max_{u_t} \{r(x_t, X_t, Z_t, u_t) + \beta V(x_{t+1}, X_{t+1}, Z_{t+1})\}$$

s.t.

$$x_{t+1} = g(x_t, X_t, Z_t, u_t) \quad (130)$$

$$X_{t+1} = G(X_t, Z_t) \quad (131)$$

$$Z_{t+1} = F(Z_t) \quad (132)$$

- The function g describes the effect of the agents control variable on his own state
- The functions G and F are the **perceived law of motion** and describe the agent's beliefs about the evolution of aggregate states

The solution

- The solution of the problem will take the form

$$u_t = h(x_t, X_t, Z_t) \quad (133)$$

- In a representative agent framework, we will have $x_t = X_t$ in the solution so that

$$X_{t+1} \equiv G_A(X_t, Z_t) = g(X_t, X_t, Z_t, h(X_t, X_t, Z_t)) \quad (134)$$

- Note: symmetry/representativeness is always imposed **after** computing the individual policy functions

Recursive competitive rational expectations equilibrium

Definition 2 (Recursive competitive rational expectations equilibrium)

A recursive competitive rational expectations equilibrium consists of a policy function h , an actual aggregate law of motion G_A , and a perceived aggregate law G such that

- i) given the perceived law G , the policy function h solves the agent's optimization problem
 - ii) the policy function h implies that $G_A = G$, i.e. the perceived and actual law of motion are consistent
- Although terminologically imprecise, shorthand names found in the literature are **recursive competitive equilibrium** or **rational expectations equilibrium**

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