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# Essential numerical tools and perturbation analysis (2.a)

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## Day 2: Solving the Model

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## The Linear Model

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Now we want to solve the linear model (we drop the hats):

$$E_t [Ay_{t+1} + By_t + Cy_{t-1} + De_t] = 0$$

where the *recursive* solution is

$$y_t = G_y y_{t-1} + G_e e_t$$

If  $y_t \in \mathbf{R}^n$  and  $e_t \in \mathbf{R}^{n_e}$  then:

- $A, B, C, G_y \in \mathbf{R}^n \times \mathbf{R}^n$
- $D, G_e \in \mathbf{R}^n \times \mathbf{R}^{n_e}$

## The Problem

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Note that if the decision rule satisfies:  $y_t = G_y y_{t-1} + G_e e_t$

we have

$$y_{t+1} = G_y y_t + G_e e_{t+1} = G_e e_{t+1} + G_y G_e e_t + G_y G_y y_{t-1}$$

and if we make all substitutions in  $E_t [Ay_{t+1} + By_t + Cy_{t-1} + De_t]$ , we get:

$$(AG_y G_y + BG_y + C)y_{t-1} + (AG_y G_e + BG_e + DG_e)e_t = 0$$

This must be true for any  $\mathbf{y}_{t-1}$  or  $\mathbf{e}_t$ . This yields the conditions that define  $\mathbf{G}_y$  and  $\mathbf{G}_e$

$$\mathbf{A}\mathbf{G}_y^2 + \mathbf{B}\mathbf{G}_y + \mathbf{C}$$

$$\mathbf{A}\mathbf{G}_y\mathbf{G}_e + \mathbf{B}\mathbf{G}_e + \mathbf{D}$$

## The Riccati Equation

The transition matrix  $\mathbf{G}_e$  must satisfy a *second order matrix equation*:

$$\mathbf{A}\mathbf{X}^2 + \mathbf{B}\mathbf{X} + \mathbf{C}$$

From our intuition in dimension 1, we know there must be multiple solutions

- how do we find them?
- how do we select the right ones?

Obviously, the qualitative dynamics of the model are given by  $\mathbf{y}_t = \mathbf{X}\mathbf{y}_{t-1}$

For the solution to the model to be stationary, the spectral radius of  $\mathbf{X}$  should be smaller than 1.

## The State-Space System

It is possible to associate a *linear* system to this Riccati equation.

It is the *state-space* representation. It characterizes vectors  $\mathbf{v}_t = (\mathbf{y}_t, \mathbf{y}_{t+1})$  along any admissible trajectory. These vectors must satisfy:

$$\underbrace{\begin{bmatrix} \mathbf{I} & \mathbf{0} \\ \mathbf{0} & \mathbf{A} \end{bmatrix}}_{\mathbf{F}} \mathbf{v}_{t+1} = \underbrace{\begin{bmatrix} \mathbf{0} & \mathbf{I} \\ -\mathbf{C} & -\mathbf{B} \end{bmatrix}}_{\mathbf{G}} \mathbf{v}_t$$

In particular, we are interested in *fundamental* trajectories, such that  $\mu \mathbf{v}_{t+1} = \lambda \mathbf{v}_t$  where  $\mu, \lambda \in \mathbf{R}$ .

### Warning:

The formulation with a pair of generalized eigenvalues  $\mu, \lambda$  is just a technicality meant to avoid infinite eigenvalues in the calculations which can happen when  $\mathbf{A}$  is defective. To build the intuition, it is suggested to look at the case  $\mu = 1$  and  $\mathbf{A} = \mathbf{I}$ .

Note that, on a fundamental trajectory, we have  $\mu(y_t, y_{t+1}) = \lambda(y_{t-1}, y_t)$ .

These trajectories are clearly recursive:  $y_t = \frac{\lambda}{\mu} y_{t-1}$

When  $\mu = 0$  and  $\lambda \neq 0$  we say there is an infinite eigenvalue. Most of the theory works if we forget about  $\mu$  but consider only  $\lambda \in [0, \infty]$

## The eigenvalues of the system

According to generalized eigenvalue theory, the system has generically  $2n$  fundamental trajectories:  $(\mu_1, \lambda_1, v_1), \dots, (\mu_{2n}, \lambda_{2n}, v_{2n})$

To simplify our reasoning we can assume that eigenvalues are ranked in increasing eigenvalues (with infinite eigenvalues last):

$$0|\lambda_1| \leq \dots \leq |\lambda_{2n}| \leq \infty$$

Remember that fundamental trajectories are recursive?

It can be shown that any recursive solution  $X$  to the quadratic system is obtained, by selecting  $n$  different eigenvectors.

As a result, there are exactly  $\binom{2n}{n}$  different solutions to our system.

The model is **well defined** when only 1 of all this solutions is non divergent. This is equivalent to say:

$$0 \leq |\lambda_1| \leq \dots \leq \lambda_n \leq 1 < |\lambda_{n+1}| \leq \dots \leq |\lambda_{2n}| \leq \infty$$

## Example 1

Forward looking inflation:

$$\pi_t = \alpha \pi_{t+1}$$

with  $\alpha > 1$ . Is it well defined?

We can rewrite the system as:  $\alpha\pi_{t+1} - \pi_t + 0\pi_{t-1} = \pi_{t+1} - \left(\frac{1}{\alpha} + 0\right)\pi_t + \left(\frac{1}{\alpha}0\right)\pi_{t-1}$

The eigenvalues are  $0 \leq 1 < \frac{1}{\alpha}$ . The unique solution is  $\pi_t = 0\pi_{t-1}$

## Example 2

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Debt accumulation equation by a rational agent:

$$b_{t+1} - \left(1 + \frac{1}{\beta}\right)b_t + \frac{1}{\beta}b_{t-1} = 0$$

Is it well-defined?

The associated polynomial  $x^2 - \left(1 + \frac{1}{\beta}\right)x + \frac{1}{\beta}$  has two eigenvalues  $\lambda_1 = 1 < \lambda_2 = \frac{1}{\beta}$

The unique solution is  $b_t = b_{t-1}$ .

- it is a *unit-root*: any initial deviation in  $b_{t-1}$  has persistent effects

## Example 3

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Productivity process:  $z_t = \rho z_{t-1}$  with  $\rho < 1$

The generalized eigenvalues are  $\lambda_1 = \rho \leq 1 < \lambda_2 = \infty$

More generally, any variable that does not appear in  $t + 1$  creates one infinite generalized eigenvalue.

### Tip

To see where the hidden eigenvalue comes from: make  $\lambda \rightarrow \infty$  in the following equation:

$$z_{t+1} - (\lambda + \rho)z_t + \frac{\lambda}{\rho}z_{t-1} = 0$$

## Blanchard-Kahn Criterium

Remember the criterium for well-definedness?

$$0 < |\lambda_1| \leq \dots \leq \lambda_n \leq 1 < |\lambda_{n+1}| \leq \dots \leq |\lambda_{2n}| \leq \infty$$

Now realize (or admit) that for each variable not appearing in  $t + 1$  in the model, there is an associated infinite eigenvalue.

We can deduce from that a common formulation of the Blanchard-Kahn criterium:

“

The model satisfies the Blanchard-Kahn criterium if the number of eigenvalues greater than one, is exactly equal to the number of variables *appearing* in  $t + 1$ .

”

It is equivalent to the **existence and unicity of a non-divergent recursive solution**.

## Computing the solution

There are several classical methods to compute the solution to the algebraic Riccati equation:

$$AX^2 + BX + C = 0$$

- qz decomposition
  - traditionnally used in the DSGE literature
  - a little bit unintuitive but easy to implement from the state-space representation
  - constructive: it produces *all* eigenvalues which makes it easy to check BK conditions

- cyclic reduction
  - more adequate for big models
- linear time iteration
  - very easy to remember/implement

## Checking the solution

Cyclic Reduction and Linear Iteration are iterative algorithms that usually converge to a solution  $\mathbf{X}$  but sometimes fail to do so.

After using one of these algorithms we can check

- that the solution is non divergent:

$$\rho(\mathbf{X}) < 1$$

- check that the first rejected eigenvalue is smaller than 1:

$$\rho((\mathbf{A}\mathbf{X} + \mathbf{B})^{-1}\mathbf{A}) < 1$$

```

1 md"""## Checking the solution
2
3 Cyclic Reduction and Linear Iteration are iterative algorithms that usually converge
  to a solution  $\mathbf{X}$  but sometimes fail to do so.
4
5 After using one of these algorithms we can check
6 - that the solution is non divergent:
7  $\rho(\mathbf{X}) < 1$ 
8 - check that the first rejected eigenvalue is smaller than 1:
9  $\rho((\mathbf{A}\mathbf{X} + \mathbf{B})^{-1}\mathbf{A}) < 1$ 
10 """

```

### Tip

Using solvant theory, it is possible to show that the eigenvalues of  $(\mathbf{A}\mathbf{X} + \mathbf{B})^{-1}\mathbf{A}$  are precisely the inverse of all the eigenvalues that have been rejected while constructing  $\mathbf{X}$

```

1 tip(md"""Using solvant theory, it is possible to show that the eigenvalues of  $(\mathbf{A}\mathbf{X} + \mathbf{B})^{-1}\mathbf{A}$ 
  are precisely the inverse of all the eigenvalues that have been rejected
  while constructing  $\mathbf{X}$ """)

```

## Linear Time Iteration (1)

Return to the Ricatti system but suppose that decision rules today and tomorrow are different:

- today:  $y_t = \bar{y} + Xy_{t-1} + G_y e_t$
- tomorrow:  $y_{t+1} = \bar{y} + \tilde{X}y_{t-1} + G_y e_t$

Then the Ricatti equation becomes:

$$A\tilde{X}X + BX + C = 0$$

## Linear Time Iteration (2)

The linear time iteration algorithm consists in solving the decision rule  $X$  today as a function of decision rule tomorrow  $\tilde{X}$ . This corresponds to the simple formula:

$$X = -(A\tilde{X} + B)^{-1}C$$

And the full algorithm can be described as:

- choose  $X_0$
- for any  $X_n$ , compute  $X_{n+1} = T(X_n) = -(AX_n + B)^{-1}C$ 
  - repeat until convergence

### Tip

Linear Time Iteration is a special case of a Bernouilli iteration

## Linear Time Iteration (3)

Starting from a random initial guess, the linear time-iteration algorithm usually converges to the solution  $X$  with the smallest modulus:

$$\underbrace{|\lambda_1| \leq \dots \leq |\lambda_n|}_{\text{Selected eigenvalues}} \leq |\lambda_{n+1}| \leq \dots \leq |\lambda_{2n}|$$

In other words, it finds the right solution when the model is well specified.



Then you just need to check that first rejected eigenvalue is greater than 1.

### Warning:

In some cases, there is no convergence. For instance if  $|\lambda_n| = |\lambda_{n+1}|$ . Or for a specific initial value  $X_0$  such that some  $AX_n + B$  is not invertible. However when the algorithm converges, it always satisfies the above condition.

## Exercise

Finish the solution of the RBC model.

**Copy and paste the code for the model from session 1.**

```
1 md"__Copy and paste the code for the model from session 1.__"
```

```
1 Enter cell code...
```

**Use ForwardDiff to compute A,B,C,D**

```
1 md"__Use ForwardDiff to compute A,B,C,D__"
```

```
1 Enter cell code...
```

**Implement the time-iteration algorithm to solve for  $G_y$**

```
1 md"__Implement the time-iteration algorithm to solve for $G_y$__"
```

```
1 Enter cell code...
```

**Check that the solution solves the original problem**

```
1 md"__Check that the solution solves the original problem__"
```

```
1 Enter cell code...
```

**Check that the greatest eigenvalue of the solution is smaller than 1**

```
1 md"__Check that the greatest eigenvalue of the solution is smaller than 1__"
```

```
1 Enter cell code...
```

**Check that the first excluded eigenvalue is greater than 1.**

```
1 md"__Check that the first excluded eigenvalue is greater than 1__"
```

**Compute  $G_e$** 

```
1 md "__Compute $G_e$__"
```

```
1 Enter cell code...
```

**Bonus: compute the generalized eigenvalues of state-space system. Are they consistent with what you have found?**

```
1 md "__Bonus: compute the generalized eigenvalues of state-space system. Are they  
consistent with what you have found?__"
```

```
1 Enter cell code...
```

**Bonus: plot some impulse response functions.**

```
1 md ""__Bonus: plot some impulse response functions.__""
```