

**Solving DSGE models using the Blanchard-Kahn algorithm  
(with extensions)**

A typical macro model with rational expectations can be written as follows:

$$A_1 \begin{bmatrix} X_{t+1} \\ E_t P_{t+1} \end{bmatrix} = A_0 \begin{bmatrix} X_t \\ P_t \end{bmatrix} + \gamma Z_{t+1} \quad (1)$$

where:

$X_t$ : vector  $n \times 1$ , collecting variables predetermined at time  $t$ , including shock processes (usually of an AR(1) type), and other backward-looking variables; these variables are often referred to as states

$P_t$ : vector  $m \times 1$ , collecting forward-looking variables, i.e. those entering the model in form of expectations; these variables are often referred to as jumpers

$Z_t$ : vector  $k \times 1$ , collecting shocks (white noise)

$A_1, A_0$ : matrices  $(n + m) \times (n + m)$

$\gamma$ : matrix  $(n + m) \times k$

Let us assume for a while that  $A_1$  is nonsingular (invertible).<sup>1</sup> Then, system (1) can be rewritten as:

$$\begin{bmatrix} X_{t+1} \\ E_t P_{t+1} \end{bmatrix} = A \begin{bmatrix} X_t \\ P_t \end{bmatrix} + R Z_{t+1} \quad (2)$$

where:  $A = A_1^{-1} A_0$  and  $R = A_1^{-1} \gamma$ .

Using the Jordan decomposition,  $A$  can be decomposed as follows:

$$A = C \Lambda C^{-1}$$

where:

$\Lambda$ : diagonal matrix with eigenvalues of  $A$  on its main diagonal, sorted with increasing absolute value; henceforth we will assume that  $A$  has  $n + m$  distinct eigenvalues

$C$ : corresponding matrix of eigenvectors

Premultiplying (2) by  $C^{-1}$  yields:

$$C^{-1} \begin{bmatrix} X_{t+1} \\ E_t P_{t+1} \end{bmatrix} = \Lambda C^{-1} \begin{bmatrix} X_t \\ P_t \end{bmatrix} + C^{-1} R Z_{t+1} \quad (3)$$

Let us now write  $\Lambda$  in block form:

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<sup>1</sup>Singularity of  $A_1$  can often be avoided by reducing the number of equations (and hence endogenous variables) so that (1) does not include static relationships (i.e. consisting of current period variables only).

$$\Lambda = \begin{bmatrix} \Lambda_1 & 0 \\ 0 & \Lambda_2 \end{bmatrix}$$

where:

$\Lambda_1$ : matrix  $n^* \times n^*$ , collecting eigenvalues lying within the unit circle (i.e. the modulus smaller or equal to unity)

$\Lambda_2$ : matrix  $m^* \times m^*$ , collecting eigenvalues outside the unit circle

Similarly for further use, let us decompose  $A$ ,  $C^{-1}$  and  $R$ :

$$A = \begin{bmatrix} A_{11[n \times n]} & A_{12[n \times m]} \\ A_{21[m \times n]} & A_{22[m \times m]} \end{bmatrix} \quad C^{-1} = \begin{bmatrix} C_{11[n^* \times n]} & C_{12[n^* \times m]} \\ C_{21[m^* \times n]} & C_{22[m^* \times m]} \end{bmatrix} \quad R = \begin{bmatrix} R_{1[n \times k]} \\ R_{2[m \times k]} \end{bmatrix}$$

Now we define transformed variables  $Y_{t[n^* \times 1]}$  and  $Q_{t[m^* \times 1]}$ :

$$\begin{bmatrix} Y_t \\ Q_t \end{bmatrix} = C^{-1} \begin{bmatrix} X_t \\ P_t \end{bmatrix} \quad (4)$$

which allows us to rewrite (3) as:

$$\tilde{E}_t \begin{bmatrix} Y_{t+1} \\ Q_{t+1} \end{bmatrix} = \Lambda \begin{bmatrix} Y_t \\ Q_t \end{bmatrix} + C^{-1} R Z_{t+1} \quad (5)$$

where  $\tilde{E}_t$  is modified expectations operator such that  $\tilde{E}_t X_{t+1} = X_{t+1}$  and  $\tilde{E}_t P_{t+1} = E_t P_{t+1}$ .

Iterating the bottom blocks of (5) forward, applying expectations operator  $E_t$  and using  $E_t Z_{t+j+1} = 0$  yields (for  $j \geq 0$ ):

$$E_t Q_{t+j+1} = \Lambda_2^j Q_t \quad (6)$$

Note that eigenvalues collected in  $\Lambda_2$  lie outside the unit circle. Then, equation (6) is nonexploding if and only if  $Q_t = 0 \forall t$ . Using (4) to go back to our original variables and using the block structure of  $C^{-1}$ , this implies:

$$C_{21} X_t + C_{22} P_t = 0 \quad (7)$$

*Theorem 1: If the following conditions are satisfied:*

(i) *the number of eigenvalues of  $A$  lying outside the unit circle is equal to the number of jumpers, i.e.  $m^* = m$  (Blanchard-Kahn condition)*

(ii)  *$C_{22}$  is of full rank (rank condition)*

*then there exist a unique solution to (2).*

The proof can be presented as follows. If both conditions of Theorem 1 hold, (7) allows us to determine uniquely the value of jumpers for any given states:

$$P_t = -C_{22}^{-1} C_{21} X_t \quad (8)$$

Moving back to (2), the law of motion for state variables can be written as:

$$X_{t+1} = (A_{11} - A_{12}C_{22}^{-1}C_{21}) X_t + R_1 Z_{t+1} \quad (9)$$

Hence, we have the full and unique solution to the original dynamic system (2). The solution has a recursive form, given by (8) and (9).

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What if the conditions of Theorem 1 are not satisfied?

*Theorem 2: If the number of eigenvalues of  $A$  lying outside the unit circle is larger than the number of jumpers, i.e.  $m^* > m$ , or  $m^* = m$  but  $C_{22}$  is rank deficient, then there exist no stable solution to (2).*

Proof: note that (7) imposes (for given  $X_t$ )  $m^*$  restrictions on  $m$  elements of  $P_t$ . Hence, for  $m^* > m$  a stable (nonexploding) solution does not exist. Similarly, if  $C_{22}$  is square but rank deficient, system (7) does not have a solution.

*Theorem 3: If the number of eigenvalues of  $A$  lying outside the unit circle is smaller than the number of jumpers, i.e.  $m^* < m$ , then there exist infinitely many solutions to (2).*

Proof: As in the case of Theorem 2, this time there exist infinitely many vectors  $P_t$  that satisfy (7) for any  $X_t$ . In particular, the realized value of  $P_t$  may depend on information outside of the model, i.e. shocks not included in  $Z_t$  (so-called sunspot shocks) may affect the solution.

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The algorithm presented above relies on transforming the original problem (1) into (2), which requires invertibility of  $A_1$ . What if this condition does not hold? The solution is to use the Schur decomposition (so-called QZ decomposition), which allows to write (1) as:

$$Q'\Lambda Z' \begin{bmatrix} X_{t+1} \\ E_t P_{t+1} \end{bmatrix} = Q'\Omega Z' \begin{bmatrix} X_t \\ P_t \end{bmatrix} + \gamma Z_{t+1} \quad (10)$$

where  $\Omega$  and  $\Lambda$  are upper triangle matrices, while  $Q$  and  $Z$  are unitary matrices ( $QQ' = I$ ,  $ZZ' = I$ ).

The ordering is now according to the generalized eigenvalues  $\frac{\omega_{ii}}{\lambda_{ii}}$ , where  $\omega_{ii}$  and  $\lambda_{ii}$  are  $i$ -th elements of the main diagonals of  $\Omega$  and  $\Lambda$ .

Following the same steps as in the original Blanchard-Kahn algorithm we obtain:

$$\tilde{E}_t \begin{bmatrix} Y_{t+1} \\ Q_{t+1} \end{bmatrix} = \Lambda^{-1}\Omega \begin{bmatrix} Y_t \\ Q_t \end{bmatrix} + \Lambda^{-1}Q\gamma Z_{t+1} \quad (11)$$

Since  $\Lambda^{-1}\Omega$  is triangular, system (11) has a recursive form and can be solved as presented before.

The algorithm using the QZ decomposition can be implemented with routine `gensys` by C. Sims, coded for Matlab/Octave and R. Documentation and codes can be downloaded from: <http://sims.princeton.edu/yftp/gensys/>.