Lecture 6 - Forecasting with VAR Models

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CENTRO DE ESTUDOS QUANTITATIVOS EM ECONOMIA E FINANÇAS

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Introduction

- A natural extension of the univariate time series models is the vector (multivariate) *ARMA* class (*VARMA*).
- The main advantage of these models is their generality and they do not impose any strong a priori restrictions on the cross-variable relationships.
- The cost is that they are heavily parameterized .
- The seminal work by [Sims, 1980] led to a widespread use of VAR models.
- Since it is difficult to model and estimate a multivariate MA component, usually only Vector Autoregressive (VAR) models will be considered.

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Representation I

• A VAR(p) model for the set of m variables y_{1t}, \dots, y_{mt} grouped in the $m \times 1$ vector $\mathbf{y}_t = (y_{1t}, \dots, y_{mt})'$ is given by

$$\mathbf{y}_{t} = \mathbf{c}_{m \times 1} + \Phi_{1} \mathbf{y}_{t-1} + \dots + \Phi_{p} \mathbf{y}_{t-p} + \epsilon_{t} \text{ and } \epsilon_{t} \sim WN\left(0, \sum_{m \times m}\right)$$
(1)

- In a VAR(p) each variable depends on:
 - up to p of its own lags and up to p lags of each other variables
 - an intercept or other deterministic components seasonal dummies or a time trend
 - error term of each equation has zero mean and is serially uncorrelated and homoskedastic but it can be contemporaneously correlated with the erros in other equations
- The number of parameters for n a VAR(p) is $\frac{m}{\text{intercepts}} + \frac{m^2p}{\text{coeficients of lagged variable}} + \frac{m(m+1)/2}{\text{var-cov for errors}}$

Representation II

- for example if m = 10 and p = 2 we have 265 parameters
- As in the univariate case, weak stationarity is an important property. It requires that

$$E(\mathbf{y}_t) = \mathbf{c}^*$$
 $Var(\mathbf{y}_t) = \Gamma_0 < \infty$
 $Cov(\mathbf{y}_t, \mathbf{y}_{t+k}) = \Gamma_k$ depends on k but not on t (2)

where $\mathbf{c}^* = (\mathbf{I} - \Phi_1 - \dots \Phi_p)^{-1} \times \mathbf{c}$.

• For a VAR(p) process, weak stationarity is verified when all the roots z of $\det(\mathbf{I} - \Phi_1 \mathbf{z} - \cdots - \Phi_p \mathbf{z}^p) = 0$ are larger than one in absolute value.

Representation III

• To theoretically justify the use of VAR models we should use the Wold representation theorem for the multivariate case that for a \mathbf{y}_t m-dimensional weakly stationary process, that is, it admits the representation

$$\mathbf{y}_t = \mathbf{C}(L)\epsilon_t \tag{3}$$

• where C(L) is a $m \times m$ matrix polynomial in the lag operator L, that is

$$\mathbf{C}(L) = \mathbf{I} + \mathbf{C}_1 L + \mathbf{C}_2 L^2 + \cdots$$

• Under mild conditions $\mathbf{C}(L)$ can be approximated by $\mathbf{A}^{-1}(L)\mathbf{B}(L)$ and (3) can be rewrite as

$$\mathbf{A}(L)\mathbf{y}_t = \mathbf{B}(L)\epsilon_t \tag{4}$$

• with $\mathbf{A}(L) = \mathbf{I} + \mathbf{A}_1 L + \cdots + \mathbf{A}_p L^p$, and $\mathbf{B}(L) = \mathbf{I} + \mathbf{B}_1 L + \cdots + \mathbf{B}_q L^q$



Representation IV

• Under slightly more stringent condition we can assume q=0 and (4) becomes a VAR(p)

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Specification of the model

- For the specification of the number of lags use multivariate versions of the information criteria.
- The most common multivariate IC are

$$AIC(j) = \ln \left| \widehat{\Sigma}_{j} \right| + \frac{2}{T} j m^{2}$$

$$BIC(j) = \ln \left| \widehat{\Sigma}_{j} \right| + \frac{\ln(T)}{T} j m^{2} \quad j = 1, \dots, p_{\text{max}}$$

• Or, in terms of the log-likelihood:

$$\begin{split} AIC(j) &= \frac{\left(-2\widehat{\ell} + 2jm^2\right)}{T} \\ BIC(j) &= \frac{\left(-2\widehat{\ell} + \ln(T)jm^2\right)}{T} \quad j = 1, \cdots, p_{\text{max}} \end{split}$$

• For a discussion of the use of these and related scalar measures to choose between alternative models in a class, see [Lütkepohl, 2007]

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Estimation

- Since the regressors are the same in each equation of the VAR, OLS equation by equation is equivalent to OLS for the system and also equivalent to GLS for the system (see my lecture notes in Time Series Econometrics available here VAR&VEC)
- Also under normality OLS and MLE are equivalent

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Diagnostic checks

- It is important to asses whether the underlying assumptions are suported by the data.
- In particular, if the errors are multivariate white noise, namely, uncorrelated and homoskedastic
- These properties can be tested using the multivariate versions of the statistics for no serial correlation and homoskedasticity and normality

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Forecasting I

• The optimal forecast in the MSFE sense for a VAR(p) is obtained as a simple extension of the formula for the univariate case:

$$\widehat{\mathbf{y}}_{T+h} = \Phi_1 \widehat{\mathbf{y}}_{T+h-1} + \dots + \Phi_p \widehat{\mathbf{y}}_{T+h-p}$$
 (5)

- where $\widehat{\mathbf{y}}_{T+h-j} = \mathbf{y}_{T+h-j}$ for $h-j \leq 0$, can possibily add an intercept \mathbf{c} or any other deterministic component
- Hence to compute the forecast for $\hat{\mathbf{y}}_{T+h}$ we calculate $\hat{\mathbf{y}}_{T+1}$, use it to obtain $\hat{\mathbf{y}}_{T+2}$ and keep iterating until we obtain $\hat{\mathbf{y}}_{T+h}$. This approach is usually defined as "iterated forecasting".

Forecasting II

• For example if \mathbf{y}_t follows a VAR(1) then

$$\begin{array}{lcl} \widehat{\mathbf{y}}_{\mathcal{T}+1} & = & \Phi_1 \mathbf{y}_{\mathcal{T}} \\ \widehat{\mathbf{y}}_{\mathcal{T}+2} & = & \Phi_1 \widehat{\mathbf{y}}_{\mathcal{T}+1} = \Phi_1^2 \mathbf{y}_{\mathcal{T}} \\ & \vdots \\ \widehat{\mathbf{y}}_{\mathcal{T}+h} & = & \Phi_1 \widehat{\mathbf{y}}_{\mathcal{T}+h-1} = \Phi_1^h \mathbf{y}_{\mathcal{T}} \end{array}$$

• In the VAR(p) context, the direct model takes the form

$$\mathbf{y}_t = \Phi_1 \mathbf{y}_{t-h} + \dots + \Phi_p \mathbf{y}_{t-h-p} + \epsilon_t$$

with forecast

$$\widetilde{\mathbf{y}}_{T+h} = \Phi_1 \mathbf{y}_T + \dots + \Phi_p \mathbf{y}_{T-p} \tag{6}$$

Forecasting III

- Under correct specification $\hat{\mathbf{y}}_{T+h}$ is more efficient than $\tilde{\mathbf{y}}_{T+h}$. However in the presence of model mis-specifications, $\tilde{\mathbf{y}}_{T+h}$ can be more robust than $\hat{\mathbf{y}}_{T+h}$.
- In the context of forecasting the notion of Granger non-causality is important. Given a *VAR* written in the form

$$\begin{bmatrix} a_{11}(L) & a_{12}(L) \\ a_{21}(L) & a_{22}(L) \end{bmatrix} \begin{bmatrix} y_t \\ x_t \end{bmatrix} = \begin{bmatrix} \epsilon_{y_t} \\ \epsilon_{x_t} \end{bmatrix}$$
 (7)

- x_t does not Granger-cause y_t if $a_{12}(L) = 0$.
- Similarly y_t does not Granger-cause x_t if $a_{21}(L) = 0$.
- We can also write the optimal forecast using the $MA(\infty)$ representation (3) in the following way

$$\widehat{\mathbf{y}}_{T+h} = \sum_{i=0}^{\infty} \mathbf{C}_{j+h} \epsilon_{T-j} \tag{8}$$

Forecasting IV

with associated forecast error

$$\mathbf{e}_{T+h} = \sum_{j=0}^{h-1} \mathbf{C}_{j+h} \epsilon_{T+h-j} \tag{9}$$

Hence the variance-covariance matrix of the forecast error is

$$V(\mathbf{e}_{T+h}) = \Sigma + \mathbf{C}_1 \Sigma \mathbf{C}_1' + \dots + \mathbf{C}_{h-1} \Sigma \mathbf{C}_{h-1}'$$
(10)

- The elements on the diagonal of $V(\mathbf{e}_{T+h})$ can be used to construct interval forecast for variables $\hat{y}_{j,T+h}$ $j=1,\cdots,m$.
- The formula for the optimal forecast in (8) can be rearranged into

$$\widehat{\mathbf{y}}_{T+h} = \mathbf{C}_h \epsilon_T + \sum_{j=0}^{3} \mathbf{C}_{j+h+1} \epsilon_{T-1-j} \\
= \widehat{\mathbf{y}}_{T+h|T-1} + \mathbf{C}_h \epsilon_{T+1} \epsilon_{T+1-j} \\
= \widehat{\mathbf{y}}_{T+h|T-1} + \mathbf{C}_h \epsilon_{T+1-j} + \mathbf{C}_h \epsilon_{T+1-j} \\
= \widehat{\mathbf{y}}_{T+h|T-1} + \mathbf{C}_h \epsilon_{T+1-j} + \mathbf{C}_h \epsilon_{T+1-j} \\
= \widehat{\mathbf{y}}_{T+h|T-1} + \mathbf{C}_h \epsilon_{T+1-j} + \mathbf{C}_$$

Forecasting V

• and using (9) gives

$$\widehat{\mathbf{y}}_{T+h} = \widehat{\mathbf{y}}_{T+h|T-1} + \mathbf{C}_h \left(\underbrace{\mathbf{y}_{T} - \widehat{\mathbf{y}}_{T|T-1}}_{\text{forecast error}} \right)$$
(12)

- The left hand side of equation (12) is the optimal forecast for period T+h made in period T and the right hand side is optimal forecast for period T+h made in period T-1 and one-step ahead forecast error made in forecasting y_T in period T-1
- It was assumed that the parameters were known
- When the parameters are estimated, the expressions for the forecast error variance must be modified to take into account parameter estimation uncertainty.
- See [Lütkepohl, 2007] for a detailed deviation of these expressions

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3.5 Forecasting with Estimated Processes

3.5.1 General Assumptions and Results

In Chapter 2, Section 2.2, we have seen that the optimal h-step forecast of the process (3.1.1) is

$$y_t(h) = \nu + A_1y_t(h-1) + \cdots + A_ny_t(h-p),$$
 (3.5.1)

where $y_t(j) = y_{t+j}$ for $j \le 0$. If the true coefficients $B = (\nu, A_1, \dots, A_p)$ are replaced by estimators $\widehat{B} = (\widehat{\nu}, \widehat{A}_1, \dots, \widehat{A}_p)$, we get a forecast

$$\hat{y}_{t}(h) = \hat{v} + \hat{A}_{1}\hat{y}_{t}(h-1) + \cdots + \hat{A}_{p}\hat{y}_{t}(h-p),$$
 (3.5.2)

where $\hat{y}_t(j) = y_{t+j}$ for $j \le 0$. Thus, the forecast error is

$$y_{t+h} - \widehat{y}_t(h) = [y_{t+h} - y_t(h)] + [y_t(h) - \widehat{y}_t(h)]$$

$$= \sum_{i=0}^{h-1} \Phi_i u_{t+h-i} + [y_t(h) - \hat{y}_t(h)], \qquad (3.5.3)$$

(3.5.4)

where the k are the coefficient matrices of the anomical MA representation of p_0 (see (2.9)). Uniform for the process p_0 , the foreast errors can be shown to have zero mean, $E[y_{0,k} = p_0(t)] = 0$, so that the contract of the process $p_0(t) = 0$, so that the process $p_0(t) = 0$ and $p_0(t) = 0$, so that the contract of the process $p_0(t) = 0$, so that the contract of the process $p_0(t) = 0$, so that the contract of the process $p_0(t) = 0$, and the process $p_0(t) = 0$, where $p_0(t) = 0$ and $p_0(t) = 0$, where $p_0(t) = 0$ is the scanned term correspond to periods $k \ge 5$, the estimation is done with in the second term correspond to periods $k \ge 5$, the estimation is done with the process $p_0(t) = 0$ and $p_0(t) = 0$.

$$\Sigma_{\widehat{y}}(h) := MSE[\widehat{y}_t(h)] = E\{[y_{t+h} - \widehat{y}_t(h)][y_{t+h} - \widehat{y}_t(h)]'\}$$

 $= \Sigma_y(h) + MSE[y_t(h) - \widehat{y}_t(h)],$

where

$$\Sigma_y(h) = \sum_{i=0}^{b-1} \Phi_i \Sigma_u \Phi'_i$$

(see [2,2,11)]. In order to evaluate the last term in (3.5.4), the distribution of the estimator \tilde{B} is needed. Because in several mode to derive the small sample distributions of the estimator \tilde{B} is needed. Because in the set of the previous sections but we have derived the asymptotic distributions of the estimators considered in the previous sections but we have derived the asymptotic distributions instead, we cannot hope for more than an asymptotic approximation to the MSE of $y_i(h) - \bar{y}_i(h)$. Such an an accordination will be derived in the following.

There are two alternative assumptions that can be made in order to facilitate the derivation of the desired result:

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(1) Only data up to the forecast origin are used for estimation.

(2) Estimation is done using a realization (time series) of a process that is independent of the process used for prediction and has the same stochastic structure (for instance, it is Gaussian and has the same first and second moments as the process used for prediction).

The first assumption is the more resultate one from a practical point of view because estimation and ferocated, gas usually based on the same data set. In that case, because the sample size is assumed to go to infinity in deriving asymptotic results, either the forecast origin has to go to infinity to or it has to be assumed that more and more data at the beginning of the sample forms and the same of the same

$$\sqrt{T}(\hat{\beta} - \beta) \xrightarrow{d} \mathcal{N}(0, \Sigma_{\hat{R}}).$$
 (3.5.5)

Samaranayake & Hasza (1988) and Basu & Sen Roy (1986) give a formal proof of the result that the MSE approximation obtained in the following remains valid under assumption (1) above.

With the foregoing assumptions it follows that, conditional on a particular realization $Y_t = (y'_1, \dots, y'_{t-p+1})'$ of the process used for prediction,

$$\sqrt{T} \left[\hat{y}_{t}(h) - y_{t}(h) | Y_{t} \right] \stackrel{d}{\rightarrow} \mathcal{N} \left(0, \frac{\partial y_{t}(h)}{\partial \beta'} \Sigma_{\hat{\beta}} \frac{\partial y_{t}(h)'}{\partial \beta} \right)$$
 (3.5.6)

because $y_t(h)$ is a differentiable function of β (see Appendix C, Proposition C.15(3)). Here T is the sample size (time series length) used for estimation. This result suggests the approximation of MSE $[\hat{y}_t(h) - y_t(h)]$ by $\Omega(h)/T$,

$$\Omega(h) = E \left[\begin{array}{c} \frac{\partial y_t(h)}{\partial A'} \Sigma_{\hat{\beta}} \frac{\partial y_t(h)'}{\partial A} \end{array} \right].$$
 (3.5.7)

In fact, for a Gaussian process
$$y_t$$
,

$$\sqrt{T} [\hat{y}_t(h) - y_t(h)] \xrightarrow{d} \mathcal{N}(0, \Omega(h)).$$
 (3.5.8)

Hence, we get an approximation

$$\Sigma_{\bar{y}}(h) = \Sigma_{y}(h) + \frac{1}{T}\Omega(h) \qquad (3.5.9)$$

for the MSE matrix of $\hat{y}_t(h)$.



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3 Estimation of Vector Autoregressive Processes

From (3.5.7) it is obvious that $\Omega(h)$ and, thus, the approximate MSE $\Sigma_n(h)$ can be reduced by using an estimator that is asymptotically more efficient than $\hat{\beta}$, if such an estimator exists. In other words, efficient estimation is of importance in order to reduce the forecast uncertainty.

3.5.2 The Approximate MSE Matrix

To derive an explicit expression for $\Omega(h)$, the derivatives $\partial y_t(h)/\partial \beta'$ are needed. They can be obtained easily by noting that

$$y_t(h) = J_1B^hZ_t$$
, (3.5.10)

where $Z_t := (1, y'_t, \dots, y'_{t-n+1})'$,

where
$$Z_1 := (1, y_1, \dots, y_{l-2p+1})^l$$
,
$$B := \begin{bmatrix} 1 & 0 & 0 & \dots & 0 & 0 \\ \nu & A_1 & A_2 & \dots & A_{p-1} & A_p \\ 0 & A_2 & 0 & \dots & 0 & 0 \\ \vdots & \vdots & A_k & 0 & \vdots & \vdots \\ 0 & 0 & \dots & I_k & \vdots & \vdots \\ 0 & 0 & \dots & I_k & \vdots & \vdots \\ 0 & 0 & \dots & I_k & \vdots & \vdots \\ 0 & 0 & \dots & I_k & \vdots & \vdots \\ 0 & 0 & \dots & 0 & \vdots & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 0 & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 0 & \vdots & \vdots \\ 0 & 0 & 0 & \dots & 0 & \vdots & \vdots \\ 0 & 0 & 0 & 0 & \dots & 0 & \vdots \\ 0 & 0 & 0 & 0 &$$

and

$$J_1 := [\underbrace{0}_{(K \times 1)} : I_K : \underbrace{0 : \cdots : 0}_{(K \times K(p-1))}] \quad [K \times (Kp+1)].$$

The relation (3.5.10) follows by induction (see Problem 3.8). Using (3.5.10),

$$\frac{\partial y_0(h)}{\partial \mathcal{G}} = \frac{\partial \text{vec}(J, \mathbf{B}^{\mathsf{L}} \mathcal{G}_i)}{\partial \mathcal{G}} = \frac{\partial \text{vec}(J, \mathbf{B}^{\mathsf{L}} \mathcal{G}_i)}{\partial \mathcal{G}}$$

$$= (Z_i^{\mathsf{L}} \otimes J_i) \begin{cases} \sum_{k \in \mathbb{N}} (\mathbf{B}^{\mathsf{L}} \mathbf{G}^{\mathsf{L}} \mathbf{G}^{\mathsf{L}}) & \frac{\partial \text{vec}(\mathbf{B}^{\mathsf{L}})}{\partial \mathcal{G}^{\mathsf{L}}} \\ (Appendix A.13, Rule (8)) \end{cases}$$

$$= (Z_i^{\mathsf{L}} \otimes J_i) \begin{cases} \sum_{k \in \mathbb{N}} (\mathbf{B}^{\mathsf{L}} \mathbf{G}^{\mathsf{L}} \mathbf{G}^{\mathsf{L}}) & \frac{\partial \text{vec}(\mathbf{B}^{\mathsf{L}})}{\partial \mathcal{G}^{\mathsf{L}}} \\ (I_{\mathcal{E} \mathcal{F}^{\mathsf{L}}} \otimes J_i) & \frac{\partial \text{vec}(\mathbf{B}^{\mathsf{L}})}{\partial \mathcal{G}^{\mathsf{L}}} \\ & = \sum_{k \in \mathbb{N}} Z_i(\mathbf{B}^{\mathsf{L}})^{k-1-i} \otimes J_i \mathbf{B}^{\mathsf{L}} J_i^{\mathsf{L}} \\ & = \sum_{k \in \mathbb{N}} Z_i(\mathbf{B}^{\mathsf{L}})^{k-1-i} \otimes \Phi_i, \qquad (3.5.11)$$

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where $\phi_i = J_1 \mathbf{B}^i J_1^i$ follows as in (2.1.17). Using the LS estimator $\widehat{\boldsymbol{\beta}}$ with asymptotic covariance matrix $\Sigma_{\widehat{\boldsymbol{\beta}}} = \varGamma^{-1} \otimes \Sigma_w$ (see Proposition 3.1), the matrix $\Omega(h)$ is seen to be

$$\begin{split} D(h) &= E \left[\frac{g_{0}(h)}{g_{0}^{2}} (\Gamma^{-1} \otimes \Sigma_{h}) \frac{g_{0}(h)^{2}}{g_{0}^{2}} \right] \\ &= \sum_{h=1}^{h-1} \sum_{j=0}^{h-1} E(Z_{j}^{j}(\mathbf{P})^{h-1-j} \Gamma^{j} \mathbf{P}^{h-j-j} Z_{j}^{j}) \otimes \phi_{h} \Sigma_{h} \phi_{j}^{j} \\ &= \sum_{j=0}^{h-1} \sum_{j=0}^{h-1} E(Z_{j}^{j}(\mathbf{P})^{h-1-j} \Gamma^{j} \mathbf{P}^{h-j-j} Z_{j}^{j}) \phi_{h} \Sigma_{h} \phi_{j}^{j} \\ &= \sum_{j=0}^{h-1} \sum_{j=0}^{h-1} E(\mathbf{P})^{h-j-j} \mathbf{P}^{h-j-j} E(Z_{j}^{j}(\mathbf{p})_{h}^{j} \Sigma_{h}^{j} \phi_{j}^{j} \\ &= \sum_{j=0}^{h-1} \sum_{j=0}^{h-1} m(\mathbf{P})^{h-j-j} \mathbf{P}^{h-j-j} F(\mathbf{p}) \delta_{h} \Sigma_{h} \phi_{j}^{j}, \end{split}$$
(3.5.12)

provided u is stable so that

$$\Gamma := \text{plim}(ZZ'/T) = E(Z_tZ'_t).$$

Here $Z:=(Z_0,\dots,Z_{T-1})$ is the $((Kp+1)\times T)$ matrix defined in (3.2.1). For example, for h=1,

$$\Omega(1) = (Kp + 1)\Sigma_u$$
.

Hence, the approximation

$$\Sigma_{g}(1) = \Sigma_{u} + \frac{Kp + 1}{T}\Sigma_{u} = \frac{T + Kp + 1}{T}\Sigma_{u} \qquad (3.5.13)$$

of the MSE matrix of the 1-step forecast with estimated coefficients is obtained. This expression shows that the contribution of the estimation variability to the forecast MSE matrix $\Sigma_0(1)$ depends on the dimension K of the process, the VAR order p, and the sample size T used for estimation. It can be quite substantial if the sample size is small or moderate. For instance, considering a three-dimensional process of order 8 which is estimated from 15 years of quarterly data (i.e., T = 52 plus 8 presample values needed for LS estimation), the 1-step forecast MSE matrix E, for known processes is inflated by a factor (T + Kp + 1)/T = 1.48. Of course, this approximation is derived from asymptotic theory so that its small sample validity is not guaranteed. We will take a closer look at this problem shortly. Obviously, the inflation factor $(T+Kp+1)/T \rightarrow 1$ for $T \rightarrow \infty$. Thus the MSE contribution due to sampling variability vanishes if the sample size gets large. This result is a consequence of estimating the VAR coefficients consistently. An expression for $\Omega(h)$ can also be derived on the basis of the mean-adjusted form of the VAR process (see Problem 3.9).

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3 Estimation of Vector Autoregressive Processes

In practice, for h > 1, it will not be possible to evaluate D(h) without knowing the AR conflictents summarized in the matrix B(h). A consistent estimator $\tilde{H}(h)$ may be obtained by replacing all unknown parameters by their LS estimators, that is, B is replaced by B which is obtained by using B for B. \mathcal{L}_b is replaced by \tilde{L}_b , \tilde{L}_b is continued by \tilde{L}_b , \tilde

The foregoing discussion is of importance in setting up interval forecasts. Assuming that y_i is Gaussian, an approximate $(1 - \alpha)100\%$ interval forecast, h periods ahead, for the k-th component $y_{k,x}$ of y_i is $y_{k,x}(h) + y_{k,y}(x,y) = \frac{\partial y_i}{\partial x_i}(h)$ (3.5.14)

$$\hat{y}_{k,t}(h) \pm z_{(\alpha/2)}\hat{\sigma}_k(h)$$
 (3.5.14)

$$\left[\widehat{y}_{k,t}(h) - z_{(\alpha/2)}\widehat{\sigma}_{k}(h), \ \widehat{y}_{k,t}(h) + z_{(\alpha/2)}\widehat{\sigma}_{k}(h)\right],$$
 (3.5.15)

where $x_{(\alpha)}$ is the upper α 100-th percentile of the standard normal distribution and $\tilde{\sigma}_k(h)$ is the square root of the k-th diagonal element of $\tilde{E}_p(h)$. Using Bonferroni's inequality, approximate joint confidence regions for a set of forecasts can be obtained just as described in Section 2.2.3 of Chapter 2.

3.5.3 An Example

To illustrate the previous results, we consider again the investment/income/ consumption example of Section 3.2.3. Using the VAR(2) model with the coefficient estimates given in (3.2.22) and

$$y_{T-1} = y_{72} = \begin{bmatrix} .02551 \\ .02434 \\ .01319 \end{bmatrix}$$
 and $y_T = y_{73} = \begin{bmatrix} .03637 \\ .00517 \\ .00599 \end{bmatrix}$

results in forecasts

$$\begin{split} \hat{y}_{T}(1) &= \hat{y} + \hat{A}_{1}y_{T} + \hat{A}_{2}y_{T-1} = \begin{bmatrix} -.011 \\ .020 \\ .022 \end{bmatrix}, \\ \hat{y}_{T}(2) &= \hat{y} + \hat{A}_{1}\hat{y}_{T}(1) + \hat{A}_{2}y_{T} = \begin{bmatrix} .011 \\ .020 \\ .020 \end{bmatrix}, \\ .015 \end{bmatrix}, \end{split}$$
(3.5.16)

and so on

The estimated forecast MSE matrix for h = 1 is

$$j(1) = \frac{T + Kp + 1}{T} \hat{\Sigma}_{w} = \frac{73 + 6 + 1}{73} \hat{\Sigma}_{w}$$

$$= \begin{bmatrix} 23.34 & .785 & 1.351 \\ .785 & 1.506 & .674 \\ 1.331 & .674 & .978 \end{bmatrix} \times 10^{-4}, \quad (3.5.17)$$



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Impulse response functions I

• A stationary VAR can be written in $MA(\infty)$ form as

$$\mathbf{y}_{t} = \Phi^{-1}(L)\epsilon_{t} = \Theta(L)\epsilon_{t} \quad \text{with} \quad \epsilon_{t} \sim WN(0, \Sigma)$$
 (13)

ullet Since Σ is positive definite, there exists a non-singular matrix P such that

$$P\Sigma P' = I \tag{14}$$

• Can rewrite (13) as

$$\mathbf{y}_{t} = \Theta(L)P^{-1}P\epsilon_{t} = \Psi(L)v_{t}$$

$$v_{t} = P\epsilon_{t}$$

$$E(v_{t}) = 0 \text{ and } E(v_{t}v'_{t}) = P\Sigma P' = I$$

$$(15)$$

• Equation (15) is the $MA(\infty)$ representation of the model

$$P\Phi(L)\mathbf{y}_t = v_t \tag{16}$$

Impulse response functions II

- which is a Structural VAR (SVAR) as there are contemporaneous relationships among the variables because of the P matrix
- ullet The orthogonal errors v_t are the structural (economic) shocks.
- How the variable under analysis will react to the structural shocks by writing the polynomial $\Psi(L)$ as

$$\Psi(L) = P^{-1} - \Theta_1 P^{-1} L - \Theta_2 P^{-1} L^2 - \cdots
= \Psi_1 - \Psi_2 L - \Psi_3 L^2$$
(17)

Impulse response functions III

• A shock is a vector with one element equal to one and all the others equal to zero, e.g.

$$v_{1,t+1} = \begin{bmatrix} 1 \\ 0 \\ 0 \\ \vdots \\ 0 \end{bmatrix}, v_{2,t+1} = \begin{bmatrix} 0 \\ 1 \\ 0 \\ \vdots \\ 0 \end{bmatrix}, \dots, v_{m,t+1} = \begin{bmatrix} 0 \\ 0 \\ 0 \\ \vdots \\ 1 \end{bmatrix}$$
 (18)

- ullet The non-zero elements can also be set to any other constant such as the standard deviation of the corresponding element of ϵ_t
- Since the system is linear, the size of the shock is irrelevant because the response is just proportional to the size of the shock.

Impulse response functions IV

• The response in period t + i of y to a shock in period t + 1 will be:

$$\frac{\partial \mathbf{y}_{t+1}}{\partial v_{t+1}} = P^{-1} = \Psi_{1}$$

$$\frac{\partial \mathbf{y}_{t+2}}{\partial v_{t+1}} = -\Psi_{2}$$

$$\frac{\partial \mathbf{y}_{t+3}}{\partial v_{t+1}} = -\Psi_{3}$$

$$\vdots$$
(19)

• where for example

$$\Psi_{1} = \begin{bmatrix} \frac{\partial y_{1,t+1}}{\partial v_{1,t+1}} & \cdots & \frac{\partial y_{1,t+1}}{\partial v_{m,t+1}} \\ \vdots & \ddots & \vdots \\ \frac{\partial y_{m,t+1}}{\partial v_{1,t+1}} & & \frac{\partial y_{m,t+1}}{\partial v_{m,t+1}} \end{bmatrix}$$

$$(20)$$

Impulse response functions V

- and the terms $\frac{\partial y_{k,t+i}}{\partial v_{j,t+1}}$ are known as the impulse response in period t+i of variable k to the shock j with $k,j=1,\ldots,m$ and $i=1,2,\cdots$
- The collection of all impulse responses is known as the impulse response function (IRF)
- The matrix P is not unique as a consequence the IRF is also not unique.
- Since Σ has m(m+1)/2 distinct elements this is the maximum number of unrestricted elements in P. So typically P is chosen to be a triangular matrix, i.e.

$$P = \begin{bmatrix} p_{11} & 0 & \dots & 0 \\ p_{21} & p_{22} & \dots & 0 \\ \vdots & & \ddots & \\ p_{m1} & p_{m2} & \dots & p_{mm} \end{bmatrix}$$

Impulse response functions VI

• where the elements p_{ij} are such that $P\Sigma P' = I$

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Forecast error variance decomposition

• The h-steps ahead forecast error can be written as

$$\mathbf{e}_{T+h} = \epsilon_{T+h} + \Theta_1 \epsilon_{T+h-1} + \cdots + \Theta_{h-1} \epsilon_{T+1}$$

so that.

$$Var(\mathbf{e}_{T+h}) = \Sigma + \Theta_1 \Sigma \Theta_1' + \dots + \Theta_{h-1} \Sigma \Theta_{h-1}'$$

which can be rewritten as

$$Var(\mathbf{e}_{T+h}) = P^{-1}P\Sigma P'(P^{-1})' + \Theta_1 P^{-1}P\Sigma P'(P^{-1})'\Theta_1' + \cdots + \Theta_{h-1}P^{-1}P\Sigma P'(P^{-1})'\Theta_{h-1}'$$

= $\Psi_1\Psi_1' + \Psi_2\Psi_2' + \cdots + \Psi_h\Psi_h'$

It follows that

$$\Psi_{ij,1}^2 + \Psi_{ij,2}^2 + \dots + \Psi_{ij,h}^2 \tag{21}$$

- \bullet represents the contribution of the innovations in the i^{th} variable in explaining the h-steps ahead forecast error variance for y_i with $i, i = 1, 2, \dots, m$
- This is called forecast error variance decomposition (**FEVD**)

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Structural VARs with long-run restrictions I

- [Blanchard and Quah, 1989] explain how to identify VAR's imposing long-run restrictions
- Consider the bivariate VAR(1) for GDP growth and unemployment defined as

$$\mathbf{y}_t = A\mathbf{y}_{t-1} + B\upsilon_t \tag{22}$$

- where v_t are the structural shocks and $B = P^{-1}$ as before.
- The MA representation is

$$\mathbf{y}_t = (I - AL)^{-1} B v_t \tag{23}$$

• the cumulative long-run response of **y** to shocks in v is

$$(I - AL)^{-1}B = \begin{bmatrix} \pi_{11} & \pi_{12} \\ \pi_{21} & \pi_{22} \end{bmatrix} \begin{bmatrix} b_{11} & b_{12} \\ b_{21} & b_{22} \end{bmatrix}$$
(24)

Structural VARs with long-run restrictions II

- The (i, j) element of the above matrix captures the long-run effects of shock j on variable i
- From the variance-covariance matrix Σ we obtain 3 parameters so in order to indentify B we need to impose one a priori restriction on its elements.
- One possibility is to impose $b_{21} = 0$ as in the Cholesky decomposition .
- Another possibility if $\mathbf{y}_t = [\operatorname{growth}_t : \operatorname{unemployment}_t]'$ then the demand shock v_{1t} has no long-run effect on growth if

$$\pi_{11}b_{11} + \pi_{12}b_{21} = 0$$

In fact this restriction implies

$$(I - AL)^{-1}B = \begin{bmatrix} 0 & * \\ * & * \end{bmatrix} \tag{25}$$

Structural VARs with long-run restrictions III

• rather than imposing $b_{21} = 0$, restriction (40) implies that b_{11} and b_{21} are linked by that linear relationship.

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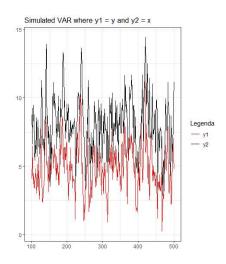
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Stationary VAR with simulated data I

- Use the following commands in R to generate a Stationary VAR
- Simulated VAR

Stationary VAR with simulated data II



Stationary VAR with simulated data III

• The VAR(1) model is given by:

$$y_t = 1 + 0.8y_{t-1} + \epsilon_{y,t}$$

$$x_t = 1 + 0.6y_{t-1} + 0.5x_{t-1} + \epsilon_{x,t}$$
(26)

with

$$v_t = \begin{bmatrix} 1 & 0.5 \\ 0 & 1 \end{bmatrix} \epsilon_t \quad \text{and} \quad v_t^{i.i.d.} N(0, I)$$
 (27)

- Then $\epsilon_t^{i.i.d.} N \left(\begin{bmatrix} 0 \\ 0 \end{bmatrix} : \begin{bmatrix} 1.25 & 0.5 \\ 0.5 & 1.0 \end{bmatrix} \right)$.
- Also note that both variables are stationary because the eigenvalues of the matrix $\begin{bmatrix} 0.8 & 0 \\ 0.6 & 0.5 \end{bmatrix}$ are $\lambda_1 = 0.8$ and $\lambda_2 = 0.5$.
- The sample size is T=600 but the estimation sample will be T=101,500 and the forecast sample from 501,...,600.

Stationary VAR with simulated data IV

- For the estimation period the best model, when BIC information criterion is used:
- selection AIC(n) HQ(n) SC(n) FPE(n)
- 1 1 1 1
- 1 2 3 4 5 6
 - AIC(n)-0.02913672-0.02089000 -0.004446732 0.001613109 0.02106193 0.03260397
 - HQ(n) -0.00485097 0.01958625 0.052220010 0.074470349 0.11010967 0.13784220
- SC(n) 0.03211594 0.08119777 0.138476141 0.185371089 0.24565502 0.29803216
- FPE(n) 0.97128426 0.97932948 0.995570936 1.001631084 1.02131634 1.03319315

Stationary VAR with simulated data V

- 7 8 9 10 11 12
- AIC(n) 0.0414267 0.0566401 0.06334653 0.07343361 0.08358958 0.08967472
- HQ(n) 0.1628554 0.1942593 0.21715626 0.24343384 0.26978031 0.29205595
- SC(n) 0.3476900 0.4037385 0.45128004 0.50220223 0.55319331 0.60011355
- FPE(n) 1.0423771 1.0583937 1.06556303 1.07642502 1.08748523 1.09420960
- ullet The estimated parameters for the best VAR(1) model are given by
- VAR(1)
- In order to compare the forecast of the best model with another model with mispecification, we also estimate a VAR(4)

Stationary VAR with simulated data VI

- VAR(4)
- Next test the specification of the VAR(1) model. We have the following results:
- Mispecification tests for VAR(1)
- And for the VAR(4)
- Mispecification tests for VAR(4)

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Stationary VAR with simulated data - Forecasting

- Use the following commands in R
- Forecasting VAR(1)
- RMSE and MAE for h-steps ahead VAR(1)

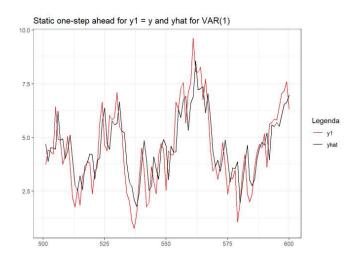
		У	X
•	RMSE	1.91	2.46
	MAE	1.60	2.07

RMSE and MAE for one-step ahead VAR(1)

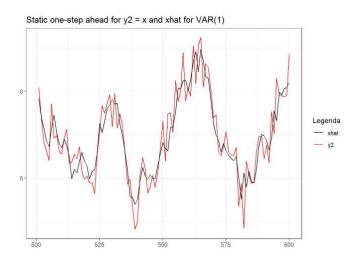
		У	X
•	$RMSE_dyn$	1.11	0.97
	$MAE_{-}dyn$	0.89	0.79

Comparison of one-step ahead and observations for VAR(1)

Stationary VAR with simulated data - Forecasting II



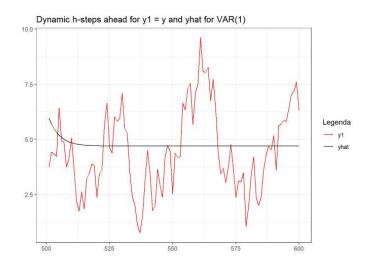
Stationary VAR with simulated data - Forecasting III



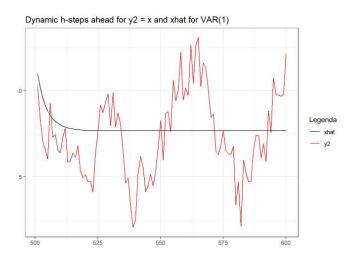
Stationary VAR with simulated data - Forecasting IV

• Comparison of h-steps ahead and observations for VAR(1)

Stationary VAR with simulated data - Forecasting V



Stationary VAR with simulated data - Forecasting VI



Stationary VAR with simulated data - Forecasting VII

- Estimate univariate ARMA models for y and x
- The following models are estimated

$$y_t = 4.9622 + u_t$$

$$u_t = 0.7684u_{t-1} + \epsilon_t$$

$$(0.0317)$$

and

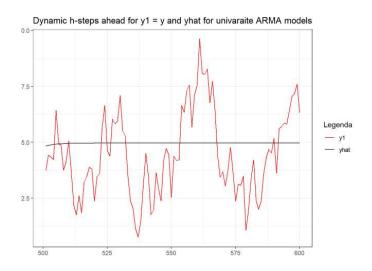
 RMSE and MAE for h-steps ahead univariate ARMA models for each variables

Stationary VAR with simulated data - Forecasting VIII

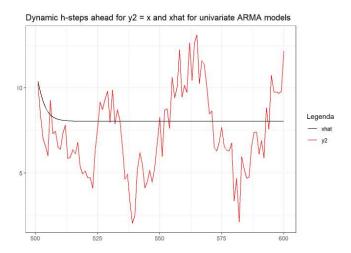
		У	×
•	RMSE	1.94	2.51
	MAE	1.64	2.11

 Comparison of h-steps ahead and observations for unvariate ARIMA for each variables

Stationary VAR with simulated data - Forecasting IX



Stationary VAR with simulated data - Forecasting X

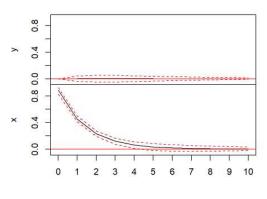


Stationary VAR with simulated data - Forecasting XI

• Impulse Response Functions (IRF) for the system (y, x)

Stationary VAR with simulated data - Forecasting XII

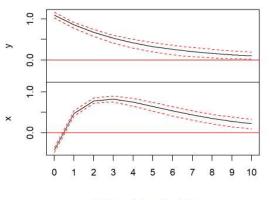
Orthogonal Impulse Response from x



95 % Bootstrap CI, 100 runs

Stationary VAR with simulated data - Forecasting XIII

Orthogonal Impulse Response from y



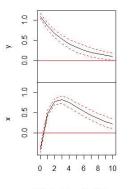
95 % Bootstrap CI, 100 runs

Stationary VAR with simulated data - Forecasting XIV

• Impulse Response Functions (IRF) for the system (x, y)

Stationary VAR with simulated data - Forecasting XV

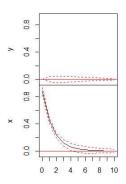




95 % Bootstrap CI, 100 runs

Stationary VAR with simulated data - Forecasting XVI





95 % Bootstrap Cl, 100 runs

Stationary VAR with simulated data - Forecasting XVII

RMSE and MAE for h-steps ahead VAR(4)

		У	X
•	RMSE4	1.92	2.47
	MAE4	1.61	2.08

• RMSE and MAE for one-step ahead VAR(4)

		У	X
•	RMSE_dyn4	1.07	0.97
	MAE_dyn4	0.88	0.78
	IVI/ (L_dyll+	0.00	0.70

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