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COMBINING INFORMATION IN STATISTICAL MODELLING

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Abstract

How to combine information from different sources is becoming an important statistical area of research under the name of Meta Analysis. This paper shows that the estimation of a parameter or the forecast of a random variable can also be seen as a process of combining information. It is shown that this approach can provide some useful insights on the robustness properties of some statistical procedures, and it also allows the comparison of statistical models within a common framework. Some general combining rules are illustrated using examples from ANOVA analysis, diagnostics in regression, time series forecasting, missing value estimation and recursive estimation using the Kalman Filter.

Key Words

Analysis of variance; Diagnostics; Forecasting; Kalman filter; Linear regression; Meta-Analysis; Time Series.

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1. INTRODUCTION

The proliferation of statistical studies in many areas of research has led to a growing interest in developing methods of combining information from different studies. This area of research was named Meta-Analysis by Glass (1976) and it has received considerable attention in the Social Sciences (Hedges and Olkin, 1985; Wolf, 1986). Examples of the use of Meta-Analysis in other scientific areas can be found in Utts (1991), Mosteller and Chalmers (1992), Dear and Begg (1992), Hedges (1992) and the references included in these papers.

The process of estimation of an unknown quantity, θ , that can be a fixed parameter or a random variable, can always be seen as a process of combining information from the data about θ . Understanding this process is crucial to evaluate the performance of an estimation rule. Often, we have independent sources of information about θ . For instance, a sample of size n can be considered as a set of j independent samples of size n_j , with $\sum n_j = n$. If we have unbiased and independent estimates of the unknown quantity, $\hat{\theta}_1, \dots, \hat{\theta}_n$ they are usually combining according to the following well known rule

Rule 1. Given n unbiased and independent estimates $\hat{\theta}_i$ of a scalar parameter θ with non zero variances σ_i^2 , the best (minimum variance) linear unbiased estimate (BLUE) of θ , $\hat{\theta}_T$, is given by

$$\hat{\theta}_T = \sum_{i=1}^n \frac{\sigma_i^{-2}}{\sum \sigma_j^{-2}} \hat{\theta}_i \quad (1.1)$$

and the variance of the pooled estimate, $\hat{\theta}_T$, is given by $(\sum \sigma_j^{-2})^{-1}$.

This rule is commonly applied in Meta-Analysis for the parametric estimation of effect size from a series of experiments (see Hedges and Olkin, Chp.6). This paper generalizes this rule for dependent and vector-valued unknown quantities and apply it to several common statistical estimation problem that are presented as particular cases of the general problem of combining different sources of information. It is shown that this approach provides some insights about the properties of the procedures considered. Also, it provides

a common ground to compare several models and estimation procedures.

The paper is organized as follows. In section 2 we show that looking at ANOVA from the perspective of rule I allows a simple understanding of the robustness properties of the estimators and of the importance of equal sample size in all groups. Section 3 shows that this approach is useful to compare two models for forecasting growth in a time series. Section 4 analyzes the estimation of missing values in linear time series and shows how this approach leads to a simple solution for dealing with the end effects. Section 5 discusses how the structure of an estimator in linear regression can suggest new diagnostics to evaluate the data robustness of the fitted model. Section 6 presents the more general rule for combining information used in the paper and applies it to derive recursive estimators and diagnostic measures. Finally, section 7 includes some concluding remarks.

2. ROBUSTNESS IN ANOVA PROBLEMS

Suppose we have two independent samples (x_1, \dots, x_n) , (y_1, \dots, y_m) from the same population and we want to estimate its mean and variance. Assuming normality, and calling \bar{x} , \bar{y} , the sample means, and s_1^2 and s_2^2 the unbiased sample variances, the application of rule I leads to

$$\hat{\mu} = \frac{n}{n+m} \bar{x} + \frac{m}{n+m} \bar{y} \quad (2.1)$$

and

$$s_T^2 = \frac{(n-1)s_1^2 + (m-1)s_2^2}{n+m-2}. \quad (2.2)$$

The result in (2.2) follows because in normal samples $\text{Var}(s^2) = 2\sigma^4/(n-1)$. When the population is not normal $\hat{\mu}$ is still the best linear unbiased estimator, whereas s_T^2 is not. This happens because the variance of \bar{x} is always σ^2/n and then rule I always leads to (2.1), whatever the parent population. However, the variance of s_T^2 for nonnormal populations is usually a more complex function of n : for instance, when the population is χ^2 , it is given by

$\sigma^4/g(n)$, where $g(n)$ is an increasing function of n . Therefore, for non normal populations the general estimate of σ^2 given by rule I is

$$s_T^2 = \frac{g(n)}{g(n) + g(m)} s_1^2 + \frac{g(m)}{g(n) + g(m)} s_2^2. \quad (2.3)$$

If $n=m$, (2.2) and (2.3) are both equal to $(s_1^2 + s_2^2)/2$, and the estimate is robust: it is BLUE whatever the population. However, if the sample sizes n and m are very different, then (2.2) and (2.3) will produce different answers.

This result will also be true in ANOVA problems. Suppose we have k different groups. Then, under the standard hypothesis of homogeneity in variance in all groups, the residual variance estimate is given by

$$s_R^2 = \sum \left[\frac{n_i - 1}{n - k} \right] s_i^2 \quad (2.4)$$

where $s_i^2 = (n_i - 1)^{-1} \sum_{j=1}^{n_i} (y_{ij} - \bar{y}_i)^2$ is the unbiased variance estimate in group i . Again, if the population is not normal (2.4) may be a very bad estimate and will be in contradiction with rule I. However, when the sample size is equal in all groups and assuming $\text{Var}(s_i^2) = \sigma^4/g(n_i)$, it will be BLUE, whatever the population, for here $g(n_i) \equiv g(n)$.

3. COMPARING ESTIMATES OF GROWTH IN TIME SERIES

Two procedures often used for forecasting the future growth of a given time series are: (i) detrend the observed data by regressing the observations on time, fit a stationary time series model to the residuals from this regression and build the forecast as the sum of the deterministic trend and the forecast of the stationary residual; (ii) difference the series, fit a stationary ARMA model in the first difference of the series and forecast the series using the ARIMA model. Typically models built in this way include a constant for many economic time series. The decision on which of these two procedures should be used is made by testing whether or not the series has one unit root. However, the available tests are not very powerful, specially for short time series, (see for instance De Jong et al 1992) and, therefore,

it is important to understand the consequences of using these models.

Let y_t be the time series data and let us assume, for the sake of simplicity, that the sample size is $n=2m+1$. Let $t=\{-m, \dots, 0, \dots, +m\}$. Then the least squares estimator of the slope in the regression on time

$$y_t = \beta_0 + \beta_1 t + u, \quad E(u) = 0, \quad \text{Var}(u) = \sigma^2 \quad (3.1)$$

is given by

$$\hat{\beta}_1 = \frac{\sum t y_t}{\sum t^2} = \left[2 \sum_{i=1}^m i^2 \right]^{-1} \sum_{i=1}^m t (y_t - y_{-t}). \quad (3.2)$$

Calling $b_t = y_t - y_{t-1}$ the observed growth at time t and after some straightforward manipulations that are shown in Peña (1995), the estimate of the slope can be written as

$$\hat{\beta}_1 = \sum_{j=1}^m \omega_j (b_j + b_{1-j}) \quad (3.3)$$

where the weights ω_j are given by

$$\omega_j = a_0 - a_1(j^2 - j) \quad j = 1, \dots, m$$

where $a_0 = 3/(2m+1)$ and $a_1 = 3/m(2m+1)(m+1)$, and add up to one. Therefore the estimated growth $\hat{\beta}_1$ is a weighted mean of all the observed growths b_j , with decreasing weight from the center of the sample. The maximum weights are given to b_1 and b_0 , that correspond to the observed growth in the middle of the sample period, and the minimum weights are given to b_m and b_{1-m} , the first and last observed growth. Note that the weight decrease quadratically from the middle of the sample.

The estimator (3.3) has an interesting interpretation. In the assumption that the linear model (3.1) holds, the $2m$ values b_t ($t = -m+1, \dots, m$) are unbiased estimates for β . The covariance matrix of these $2m$ estimates is the Toeplitz matrix:

$$V = \begin{bmatrix} 2\sigma^2 & -\sigma^2 & 0 & \dots & 0 \\ -\sigma^2 & 2\sigma^2 & -\sigma^2 & & \\ \cdot & \cdot & \cdot & \cdot & \\ \cdot & \cdot & \cdot & \cdot & -\sigma^2 \\ 0 & \dots & \dots & -\sigma^2 & 2\sigma^2 \end{bmatrix}. \quad (3.4)$$

Now we can set the following rule (see, for instance Newbold and Granger, 1974).

Rule II: given a vector $\hat{\theta}$ of unbiased estimators of a parameter θ with covariance matrix V , the best (in the mean squared sense) linear unbiased estimator of θ is given by

$$\hat{\theta}_T = (1'V^{-1}1)^{-1}(1'V^{-1}\hat{\theta}) \quad (3.5)$$

where $1' = (1 \ 1 \ \dots \ 1)$, and the variance of $\hat{\theta}_T$ is given by

$$\text{Var}(\hat{\theta}_T) = (1'V^{-1}1)^{-1} \quad (3.6)$$

This Rule II is a particular case of the Rule V that is proved in the appendix.

The inverse of the Toeplitz matrix (3.4) has been studied by Shaman (1969) who obtained the exact inverse of a first order moving average process. As V can be interpreted as the covariance matrix of a non-invertible ($\theta=1$) first order moving average process, then $V^{-1} = \{v_{ij}\}$, is given by

$$v_{ij} = \frac{i(2m-j+1)}{2n+1} \quad j \geq i, \quad i=1, \dots, 2m,$$

and $v_{ij} = v_{ji}$. Therefore

$$V^{-1} = \frac{1}{\sigma^2(2m+1)} \begin{bmatrix} 2m & 2m-1 & 2m-2 & \dots & 1 \\ 2m-1 & 2(2m-1) & 2(2m-2) & \dots & 2 \\ 2m-2 & 2(2m-2) & 3(2m-2) & \dots & 3 \\ \vdots & \vdots & \vdots & & \\ 2 & 4 & 6 & \dots & 2m-1 \\ 1 & 2 & 3 & \dots & 2m \end{bmatrix}$$

It is easy to show that the estimator (3.3) can also be obtained by applying Rule II to the

unbiased but correlated estimates b_t .

When an ARMA model is fitted to the residuals of the regression model, the equation for the h steps ahead forecast where we call $\hat{y}_t(h) = E[y_{t+h} | y_t, y_{t-1}, \dots]$ is

$$\hat{y}_t(h) = \hat{\beta}_0 + \hat{\beta}_1 h + \hat{\eta}_t(h) \quad (3.7)$$

where $\hat{\eta}_t(h)$ is the forecast of the zero mean stationary process fitted to the residuals. As for a stationary process the long run forecast converges to the mean, $\hat{\eta}_t(h) \rightarrow 0$, and the parameter $\hat{\beta}_1$ is the long-run estimated growth of the time series.

Let us compare (3.7), with the growth estimate provided by the integrated ARIMA model

$$\nabla y_t = \beta + n_t \quad (3.8)$$

where $\nabla = 1-B$ and $By_t = y_{t-1}$ and n_t follows a zero mean stationary ARMA model. Letting V denote the covariance matrix of n_t , the estimate of β in (3.8) is given by the generalized least squares estimator

$$\hat{\beta} = (1'V^{-1}1)^{-1} (1'V^{-1}b) \quad (3.9)$$

where the vector b has components $b_t = y_t - y_{t-1}$. Assuming that n_t is stationary and invertible it is well known (see Fuller 1976) that $\bar{b} = (1/(n-1)) \sum b_t$ is asymptotically unbiased for β . When n is large, the expected forecast h periods ahead is given by

$$\hat{y}_t(h) = \bar{b} h + \hat{\eta}_t(h) \quad (3.10)$$

where $\hat{\eta}_t(h)$ is the h -step ahead forecast of the stationary process n_t . As for h large the $\hat{\eta}_t(h)$ will go to zero, the long-run growth will be estimated by a weighted average with uniform weighing of the observed growths b_t .

In summary, the two models forecast future growth by using a weighted average of the observed growths in the sample. Linear regression gives minimum weight to the last observed growth and maximum weight to the center of the sample period. The ARIMA model gives uniform weighting in all the years in the sample. A comparison of the forecasting performance of these and other models used for forecasting growth can be found in Peña (1995).

4. ESTIMATING MISSING VALUES IN TIME SERIES

Suppose a Gaussian stationary time series y_t that follows the general representation

$$y_t = \sum_{i=1}^{\infty} \pi_i y_{t-i} + a_t \quad (4.1)$$

where a_t is a white noise process with variance σ_a^2 . Then, if the value y_T is missing, we can obtain an unbiased estimate of it by using

$$\hat{y}_T^{(0)} = \sum_{i=1}^{\infty} \pi_i y_{T-i} \quad (4.2)$$

and this estimate will have variance σ_a^2 . Also, from (4.1) we can write

$$y_T = \pi_j^{-1} \left(y_{T+j} - \sum_{i=1}^{\infty} \pi_i y_{T+j-i} \right) + a_T / \pi_j \quad (4.3)$$

Thus we can obtain additional unbiased estimates of y_T from (4.3) by

$$\hat{y}_T^{(j)} = \pi_j^{-1} \left(y_{T+j} - \sum_{i \neq j} \pi_i y_{T+j-i} \right) \quad (4.4)$$

with variance σ^2 / π_j^2 . As all these estimates are unbiased and independent given the observed data, the best linear unbiased estimate of the missing value y_T is readily obtained by applying Rule 1

$$\hat{y}_T = \sum_{j=0}^{\infty} \frac{\pi_j^2}{\sum \pi_j^2} \hat{y}_T^{(j)} \quad (4.5)$$

where $\pi_0 = -1$. It is easy to show (Maravall and Peña, 1995) that this estimate is equivalent to the well known expression for the missing value estimation in a gaussian stationary time series

$$\hat{y} = -\sum_{i=0}^{\infty} \rho_i^D (y_{T-i} + y_{T+i}) \quad (4.6)$$

(See Grenander and Rosenblatt, 1957, and Peña and Maravall, 1991). However, the advantage of formulation (4.5) is that it provides a clear understanding of how to proceed when the missing value is near the extremes of the series so that the two side symmetric filter (4.6) has to be truncated. Then, we have to combine (4.2) with the n-T estimates (4.4) that are available and the exact formula for the finite sample interpolator is

$$\hat{y}_{T,F} = \sum_{j=0}^{n-T} \frac{\pi_j^2}{\sum_0^n \pi_j^2} \hat{y}_T^{(j)} \quad (4.7)$$

This idea can be easily extended to groups of missing observations. We will illustrate it here with an example: suppose we have an AR(1) process in which the values y_T and y_{T+1} are missing. Then, for y_T we have the two estimates:

$$\hat{y}_T^{(0)} = \phi y_{T-1} \quad (4.7)$$

with variance σ_a^2 , and

$$\hat{y}_T^{(2)} = \phi^{-2} y_{T+2} \quad (4.8)$$

with variance $\sigma_a^2(1+\phi^2)/\phi^4$. The best linear unbiased estimate will be

$$\hat{y}_T = \frac{\phi(1+\phi^2)}{1+\phi^2+\phi^4} y_{T-1} + \frac{\phi^2}{1+\phi^2+\phi^4} y_{T+2} \quad (4.9)$$

that agrees with the general formula obtained by a different approach in Peña and Maravall (1991). The estimate of \hat{y}_{T+1} will be similar to (4.9) but with the roles of y_{T-1} and y_{T+2} reversed.

5. SENSITIVITY ANALYSIS IN REGRESSION

It is well known that in the linear regression model

$$y = \beta_0 + \beta_1 x + u, \quad E(u) = 0, \quad \text{Var}(u) = \sigma^2 \quad (5.1)$$

the least square estimate of the slope is given by

$$\hat{\beta} = \Sigma w_i b_i \quad (5.2)$$

where $w_i = (x_i - \bar{x})^2 / \Sigma (x_i - \bar{x})^2$ is a set of weights ($w_i \geq 0$, $\Sigma w_i = 1$) and the b_i are estimates of the slope that can be built up by using the sample data:

$$b_i = \frac{y_i - \bar{Y}}{x_i - \bar{x}}. \quad (5.3)$$

These estimates are not independent, because $a_x' b = 0$, where $a_x' = ((x_1 - \bar{x}) \dots (x_n - \bar{x}))$ and $b = (b_1, \dots, b_n)$. They have a singular covariance matrix

$$S_b = D_x^{-1} (I - 1/n \mathbf{1} \mathbf{1}') D_x^{-1} \sigma^2 \quad (5.4)$$

where D_x is a diagonal matrix such that the i th diagonal element is the i th element of a_x , that is $\text{diag}(D_x) = a_x$. Then, we can use the following rule.

Rule III: Given n dependent estimates $\hat{\theta}_i$ with singular covariance matrix S_θ , the best linear unbiased estimator of θ is given by

$$\hat{\theta}_T = (1' S_\theta^{-1} 1)^{-1} 1' S_\theta^{-1} \hat{\theta} \quad (5.5)$$

where S_θ^{-1} is a generalized inverse of S_θ , and the variance of the pooled estimator $\hat{\theta}_T$ is

$$\text{Var}(\hat{\theta}_T) = (1' S_\theta^{-1} 1)^{-1}.$$

It is straightforward to check that a generalized inverse of (5.4) is given by

$$S_b^- = D_x D_x \sigma^{-2}, \quad (5.6)$$

because $1'D_x = 0$, and if we apply (5.5) to (5.6) and (5.3) as $1'S_b^- 1' = 1/\sigma^2 \sum (x_i - \bar{x})^2$ we obtain (5.2). In summary (5.2) is again the BLUE estimate given the estimates b_i . This estimate can also be written as a weighted function of the estimates

$$b_{ij} = \frac{Y_i - Y_j}{x_i - x_j} \quad (5.7)$$

that are independent, and have variance $2\sigma^2/(x_i - x_j)^2$. Therefore, the BLUE based on b_{ij} must be

$$\hat{\beta} = \sum_i \sum_j \frac{(x_i - x_j)^2}{\sum_i \sum_j (x_i - x_j)^2} b_{ij} \quad (5.8)$$

and it is straightforward to show that this estimate is equivalent to (5.2).

Equations (5.2) shows that the leverage $(x_i - \bar{x})^2 / \sum (x_i - \bar{x})^2$ determines the potential influential of an observation on the estimated slope of the regression line, whereas the observed effect depends also on b_i . Since $\hat{\beta}$ is the sum of n components $w_i b_i$, the relative importance of a point in determining $\hat{\beta}$ can be measured by

$$\delta_i = n \left| \frac{w_i b_i}{\hat{\beta}} \right| = \left| \frac{(x_i - \bar{x})(y_i - \bar{y})}{S_{xy}} \right| \quad (5.9)$$

where $S_{xy} = \sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})$. Note that δ_i is a measure of the influence of a point (x_i, y_i) on the slope, whereas the usual statistics of influence, as the one due to Cook (1977), tries to identify both outliers and influential points. Also, the Cook's statistic can be very affected by masking (see Peña and Yohai, 1995) whereas δ_i is not. For instance, table 1 presents a set of artificial data with three large influential observations that are not identified neither by D_i (Cook's statistics) nor by the studentized residual (t_i) as extremes, but they are clearly indicated as the most influential on the slope by the statistic (5.9)

x	1	2	3	4	5	6	7	8	9	10	17	17	17
y	2	3	4	5	6	7	8	9	10	11	25	25	25
D _i	.25	.10	.03	.00	.00	.01	.02	.05	.08	.14	.16	.16	.16
t _i	1.4	1.0	.6	.3	-0.1	-0.4	-0.7	-1	-1.4	-1.7	.9	.9	.9
δ	1.3	1	.7	.5	.3	.1	.0	.0	.0	.0	2.6	2.6	2.6

Table 1

Consider now the multiple regression model

$$Y = X \beta + U \quad (5.10)$$

where X is nxp and we suppose to simplify the presentation and without loss of generality that all the variables have zero mean. Then, it is well known that each of the components of $\hat{\beta}$ can be written as

$$\hat{\beta}_j = \sum_{i=1}^n w_{ij} b_{ij} \quad (5.11)$$

where

$$b_{ij} = \frac{y_i - \bar{y}}{e_{ij.R}} \quad (5.12)$$

and $e_{ij.R}$ is the i th component of the vector of residuals $e_{j.R}$ obtained by regressing x_j on all the other explanatory variables. That is, if $X_{[j]}$ is a matrix without the j th column, x_j , and $\hat{\gamma} = (X'_{[j]} X_{[j]})^{-1} X'_{[j]} x_j$ is the least square estimate of this regression, then $e_j = x_j - X_{[j]} \hat{\gamma}$. The weight w_{ij} is given by $e_{ij.R}^2 / \sum e_{ij.R}^2$.

Suppose that we are mainly interested in some regression coefficient $\hat{\beta}_j$ that is of special interest. Then, the usual diagnostic statistics that look at the change on the whole vector of parameter estimates may not be useful. However, the weights w_{ij} provide a natural and simple way of looking at the potential effect of an observation. These weights can be computed from

$$w_{ij} = \frac{(x_{ij} - x_{i(\cdot)}) (X'_{(j)} X_{(j)})^{-1} X'_{(j)} x_j)^2}{x_j' (I - H_{(j)}) x_j} \quad (5.13)$$

where $H_{(j)}$ is the hat or projection matrix built without variable X_j . A plot of the variables w_{ij} can be useful to judge about the robustness of one estimate to the given sample.

As in the simple regression case a measure of the influence of point (x_i, y_i) on the estimation of $\hat{\beta}_j$ can be built by

$$\delta_i = n \frac{|w_{ij} b_{ij}|}{|\hat{\beta}_j|}$$

A different problem occurs when we have a sample of n_i data points (X_i, Y_i) , $i=1,2,\dots, n_i$ in which we have obtained $\hat{\beta}_i = (X_i' X_i)^{-1} X_i' Y_i$ with covariance $s_i^2 (X_i' X_i)^{-1}$, and we want to combine both estimates to obtain the BLUE. Then we can use the following rule.

Rule IV: If $\hat{\theta}_1$ is an unbiased estimator of θ with covariance matrix V_1 and $\hat{\theta}_2$ is also unbiased for θ with covariance V_2 and these two estimates are independent, the best linear unbiased estimator (minimizing the trace of the variance covariance matrix) is given by

$$\hat{\theta}_T = (V_1^{-1} + V_2^{-1})^{-1} V_1^{-1} \hat{\theta}_1 + (V_1^{-1} + V_2^{-1})^{-1} V_2^{-1} \hat{\theta}_2 \quad (5.15)$$

and the covariance matrix of the pooled estimator is

$$V_T^{-1} = V_1^{-1} + V_2^{-1}. \quad (5.16)$$

This rule is a particular case of Rule V that will be proved in the appendix, and generalizes rule I to the vector case. For instance, the BLUE estimate of $\hat{\beta}$ when combining two independent samples with the same parameter $\hat{\beta}$ but different residual variance is given by

$$\hat{\beta} = (X_1' X_1 / s_1^2 + X_2' X_2 / s_2^2)^{-1} (X_1' Y_1 / s_1^2 + X_2' Y_2 / s_2^2).$$

6. RECURSIVE ESTIMATION

Suppose we have a parametric model $y_t = f(x_t, \theta, a_t)$, that relates a vector of responses to a set of explanatory variables, x_t , a vector θ of p parameters and a set of unobserved random variables a_t . We will say that a *basic estimate* of θ is an estimate obtained from a sample of size p , while an *elemental estimate* of θ is an estimate obtained from a sample of size one. We will say that an estimate is proper if it is obtained from a sample of at least size p . For instance, if $\theta = (\mu, \sigma)$ and $y_t = \mu + \sigma a_t$, where a_t is a zero mean and unit variance scalar variable, the basic estimate requires $n=2$, and the elemental estimates, from a sample of size one y_1 , are given by $\hat{\mu} = y_1$, $\hat{\sigma}^2 = 0$, with a singular variance covariance matrix. In the standard regression model where β is $p \times 1$, the basic estimate of $\theta = (\beta, \sigma^2)$ requires $p+1$ data. The elemental estimate of β given a sample (y_1, x_1) of size one is obtained from $x_1' \hat{\beta}_1 = y_1$. Using the Moore-Peurose generalized inverse (see Guttman (1982)) and calling A^- to the generalized inverse of A the solution of this equation can be written as

$$\hat{\beta}_1 = (x_1')^- y_1 = (x_1' x_1)^{-1} x_1 y_1 \quad (6.1)$$

where x_1 is a $p \times 1$ column vector and will have a singular covariance matrix.

Sometimes we need to combine a proper and an elemental estimate of θ . For instance, in regression recursive estimation where we have an estimate $\hat{\beta}_{(n)}$ of (5.10) based on n data points, we observe y_{n+1} and need to revise $\hat{\beta}_{(n)}$ to obtain $\hat{\beta}_{(n+1)}$. In general, given a $p \times 1$ vector of parameters θ we will say that $\hat{\theta}_i$ is an elemental unbiased estimator of θ if (1) the covariance matrix of $\hat{\theta}_i$, V_i , is such that $\text{rank}(V_i) = 1$; (2) given p independent estimates $\hat{\theta}_i$ with covariance matrices V_i , the matrix $V_1^- + \dots + V_p^-$, where V_i^- is a generalized inverse of V_i is nonsingular; (3) Combining these p estimates by

$$\hat{\theta}_T = \sum_{i=1}^p \left(\sum_{i=1}^p V_i^- \right)^{-1} V_i^- \hat{\theta}_i \quad (6.2)$$

we obtain a basic unbiased estimator of θ . For instance, in linear regression the estimate (6.1) is elemental unbiased, because (1) the $p \times p$ covariance matrix of the estimate $\hat{\beta}_1$, V_1 , is

$V_i = x_i x_i' (x_i' x_i)^{-2} \sigma^2$ has rank equal to one; (2) $V_i^{-1} = x_i x_i' / \sigma^2$ and

$$\left| \sum_{i=1}^p V_i^{-1} \right| = \left(\frac{1}{\sigma^2} \right) \left| \sum_{i=1}^p x_i x_i' \right| = \left(\frac{1}{\sigma^2} \right) |X'X| \neq 0$$

and (3) combining $\hat{\beta}_i$ by

$$\hat{\beta}_T = \sum_{i=1}^p (\sum x_i x_i')^{-1} x_i y_i = (X'X)^{-1} X'Y \quad (6.3)$$

we obtain the basic BLUE estimate. We can generalize (6.1) as follows:

Rule V: Given n independent estimates $\hat{\theta}_i$ unbiased or elemental unbiased with covariance matrices V_i , that may be singular, the best (minimizing the trace of the covariance matrix) unbiased estimate is given by

$$\hat{\theta}_T = \sum_{i=1}^n \left(\sum_{j=1}^n V_j \right)^{-1} V_i^{-1} \hat{\theta}_i \quad (6.4)$$

where V_i^{-1} is the Moore-Penrose generalized inverse of V_i , and where we have assumed that $\sum V_i^{-1}$ is non singular. The covariance matrix of $\hat{\theta}_T$ is then easily seen to be

$$V_T^{-1} = \sum_{i=1}^n V_i^{-1}. \quad (6.5)$$

This Rule is proved in the appendix.

The application of this Rule V to recursive estimation leads directly to the Kalman Filter. To show this, let us consider the standard state space formulation of a dynamic model with observation equation

$$y_t = A_t \theta_t + \epsilon_t \quad (6.6)$$

and state equation

$$\theta_t = \Omega_t \theta_{t-1} + u_t \quad (6.7)$$

where y_t is $rx1$, A_t is rxp with $\text{rank}(A_t)=r$, ϵ_t is $N_r(0, C_t)$, Ω_t is pxp and $u_t \sim N_p(0, R_t)$. In this model, at any time t we may consider two independent estimates of θ . The first is the forecast of the state that comes from (6.7)

$$\hat{\theta}_t^{(1)} = \Omega_t \hat{\theta}_{t-1} \quad (6.8)$$

and whose covariance matrix can be obtained from

$$\theta_t - \hat{\theta}_t^{(1)} = \Omega_t (\theta_{t-1} - \hat{\theta}_{t-1}) + u_t \quad (6.9)$$

calling $I_t = \{y_1, \dots, y_t\}$ the information until time t , defining

$$V_{t-1} = E[(\hat{\theta}_t - \theta)(\hat{\theta}_t - \theta)' | I_{t-1}] \quad (6.10)$$

and letting $V_t = V_{t|t}$, we have from (6.9) that the covariance matrix of (6.8) is given by

$$V_{t-1} = \Omega_t V_{t-1} \Omega_t' + R_t. \quad (6.11)$$

The second estimate of θ at time t is obtained from (6.6) when y_t is observed. Assuming $p > r$ and $A_t A_t'$ non singular, this estimate is given by

$$\hat{\theta}_t^{(2)} = A_t' (A_t A_t')^{-1} y_t. \quad (6.12)$$

Using (6.6), it can be written

$$\hat{\theta}_t^{(2)} = A_t' (A_t A_t')^{-1} A_t \theta_t + A_t' (A_t A_t')^{-1} \epsilon_t \quad (6.13)$$

which shows that it is not unbiased for θ_t . However, it is easy to see that it is elemental unbiased, with singular covariance matrix

$$V_t^{(2)} = A_t' (A_t A_t')^{-1} C_t (A_t A_t')^{-1} A_t. \quad (6.14)$$

This matrix has a generalized inverse

$$(V_t^{(2)})^- = A_t' C_t^{-1} A_t.$$

Therefore, following rule V, the BLUE estimate will have a pooled covariance matrix

$$V_t^{-1} = V_{v_{t-1}}^{-1} + A_t' C_t^{-1} A_t \quad (6.15)$$

and the estimate will be given by

$$\hat{\theta}_T = (I - V_t A_t' C_t^{-1} A_t) \Omega_t \hat{\theta}_{t-1} + V_t A_t' C_t^{-1} y_t \quad (6.16)$$

or, as it is normally written,

$$\hat{\theta}_T = \Omega_t \hat{\theta}_{t-1} + V_t A_t' C_t^{-1} (y_t - A_t \Omega_t \hat{\theta}_{t-1}) \quad (6.17)$$

Equations (6.15) and (6.17) constitute the Kalman Filter, that appears as a particular case of Rule V.

It is interesting to stress that equation (6.7) provides a clear ground for building influence measures of the last observed data in recursive estimation. Calling $\hat{\theta}_{v_{t-1}} = \Omega_t \hat{\theta}_{t-1}$ to the forecast of θ_t with information until y_{t-1} , the change on the parameter vector due to observing y_t is given by

$$\hat{\theta}_t - \hat{\theta}_{v_{t-1}} = V_t A_t' C_t^{-1} e_{v_{t-1}} \quad (6.18)$$

where $e_{v_{t-1}} = y_t - A_t \hat{\theta}_{v_{t-1}}$ is the predicted residual. The Mahalanobis change on $\hat{\theta}_t$ will be given by

$$D_t = (\hat{\theta}_t - \hat{\theta}_{v_{t-1}})' V_t^{-1} (\hat{\theta}_t - \hat{\theta}_{v_{t-1}}) \quad (6.19)$$

that can be written as

$$D_t = e_{v,t-1}' C_t^{-1} A_t V_t A_t' C_t^{-1} e_{v,t-1}. \quad (6.20)$$

This diagnostic measure can be built for any statistical model in the state space form (6.6), (6.7) and estimated with the Kalman filter. It is straightforward to show that for regression models this statistic for the last observed point is equivalent to the one introduced by Cook (1977), whereas in ARIMA and transfer function models it is equivalent to the statistic introduced by Peña (1990, 1991).

7. CONCLUDING REMARKS

Any estimation or forecasting procedure can be seen as a way to combine the available information. In Bayesian statistics the prior information is combined with the posterior using Bayes' Theorem. In classical statistics the different pieces of sample information are weighted to obtain the final estimate. When we have unbiased estimators (or elemental unbiased) they are lineary combined to obtain the best linear unbiased estimate using as weights the (generalized) inverse covariance matrices. We have shown that analyzing estimates from this point of view can provide some useful insights on the properties of some statistical procedures.

APPENDIX

To prove Rule V, let us consider the class of unbiased estimators

$$\hat{\theta}_T = \sum_{i=1}^{n-1} A_i \hat{\theta}_i + (I - \sum_{i=1}^{n-1} A_i) \hat{\theta}_n \quad (\text{A.1})$$

such that if the $\hat{\theta}_i$ ($i = 1, \dots, n$) are unbiased, $\hat{\theta}_T$ will also be unbiased. The covariance matrix of $\hat{\theta}_T$ is

$$S_T = \sum_{i=1}^{n-1} A_i V_i A_i' + V_n - \sum_{i=1}^{n-1} A_i V_n - \sum_{i=1}^n V_n A_i' + \sum_{i=1}^n \sum_{j=1}^n A_i V_n A_j'$$

and the trace of this matrix is

$$\begin{aligned} m = \text{tr}(S_T) &= \sum_{i=1}^n \text{tr}(A_i V_i A_i') + \text{tr}(V_n) - 2 \sum_{i=1}^n \text{tr}(A_i V_n) + \\ &\sum_{i=1}^{n-1} \sum_{j=1}^{n-1} \text{tr}(A_i V_n A_j'). \end{aligned}$$

Now, if V is symmetric we have that $\partial \text{tr}(A V) / \partial A = V$, $\partial \text{tr}(A V A') / \partial A = 2 A V$, $\partial \text{tr}(A' V A) / \partial A = 2 V A$, $\partial \text{tr}(A V B) / \partial A = B' V$ and $\partial \text{tr}(B V A) / \partial A = V B'$, we have

$$\frac{\partial m}{\partial A_i} = 2 A_i V_i - 2 V_n + 2 \sum_{j=1}^{n-1} A_j V_n = 0$$

and so,

$$\begin{aligned} A_i V_i &= (I - \sum_{j=1}^{n-1} A_j) V_n \\ A_i &= (I - \sum_{j=1}^{n-1} A_j) V_n V_i^{-1} \end{aligned} \quad (\text{A.2})$$

Adding the $n-1$ equations (A.2), we obtain

$$\begin{aligned} \sum_{i=1}^{n-1} A_i &= V_n \sum_{i=1}^{n-1} V_i^{-1} - \sum_{j=1}^{n-1} A_j V_n \sum_{i=1}^{n-1} V_i^{-1}, \\ \sum_{i=1}^{n-1} A_i (I + V_n \sum_{i=1}^{n-1} V_i^{-1}) &= V_n \sum_{i=1}^{n-1} V_i^{-1} \end{aligned}$$

$$\sum_{i=1}^{n-1} A_i V_n \left(\sum_{i=1}^n V_i^{-1} \right) = V_n \sum_{i=1}^{n-1} V_i^{-1}$$

$$\sum_{i=1}^{n-1} A_i V_n = V_n \left(\sum_{i=1}^{n-1} V_i^{-1} \right) \left(\sum_{i=1}^n V_i^{-1} \right)^{-1}$$

and, inserting this result in (A.2)

$$A_i = V_n \left(I - \left(\sum_{i=1}^{n-1} V_i^{-1} \right) \left(\sum_{i=1}^n V_i^{-1} \right)^{-1} \right) V_i^{-1}$$

$$A_i = V_n \left(\sum_{i=1}^n V_i^{-1} - \sum_{i=1}^{n-1} V_i^{-1} \right) \left(\sum_{i=1}^n V_i^{-1} \right)^{-1} V_i^{-1}$$

$$A_i = \left(\sum_{i=1}^n V_i^{-1} \right)^{-1} V_i^{-1}.$$

We have assumed in the proof that all the inverse matrix involved exist; the proof is similar when some of these matrices are singular by replacing the inverse by the generalized inverse of the matrix.

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