# WATER RESOURCES research center

Publication No. 67

ESTIMATING MISSING VALUES IN MONTHLY RAINFALL SERIES

By

EFI FOUFOULA-GEORGIOU

A Thesis Presented to the Graduate Council of the University of Florida in Partial Fulfillment of the Requirements for the Degree of Master of Engineering

> University of Florida Gainesville



# **UNIVERSITY OF FLORIDA**

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# FLORIDA WATER RESOURCES RESEARCH CENTER

Research Project Technical Completion Report

Sponsored by

South Florida Water Management District

# A THESIS PRESENTED TO THE GRADUATE COUNCIL OF THE UNIVERSITY OF FLORIDA IN PARTIAL FULFILLMENT OF THE REQUIREMENTS FOR THE DEGREE OF MASTER OF ENGINEERING

# UNIVERSITY OF FLORIDA

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# Abstract of Thesis Presented to the Graduate Council of the University of Florida in Partial Fulfillment of the Requirements for the Degree of Master of Engineering

### ESTIMATION OF MISSING OBSERVATIONS IN MONTHLY RAINFALL SERIES

By

#### Efstathia Foufoula-Georgiou

December, 1982

Chairman: Wayne C. Huber Cochairman: James P. Heaney Major Department: Environmental Engineering Sciences

This study compares and evaluates different methods for the estimation of missing observations in monthly rainfall series. The estimation methods studied reflect three basic ideas:

- the use of regional-statistical information in four simple techniques:
  - mean value method (MV),
  - reciprocal distance method (RD),
  - normal ratio method (NR),
  - modified weighted average method (MWA);
- (2) the use of a univariate autoregressive moving average (ARMA) model which describes the time correlation of the series;

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(3) the use of a multivariate ARMA model which describes the time and space correlation of the series.

An algorithm for the recursive estimation of the missing values in a series by a parallel updating of the univariate or multivariate ARMA model is proposed and demonstrated. All methods are illustrated in a case study using 55 years of monthly rainfall data from four south Florida stations.

Chairman

# CHAPTER 1

# INTRODUCTION

#### Rainfall Records

Rainfall is the source component of the hydrologic cycle. As such it regulates water availability and thus land use, agricultural and urban expansion, maintenance of environmental quality and even population growth and human habitation. As Hamrick (1972) points out, water may be transported for considerable distances from where it fell as rain and may be stored for long periods of time, but with very few exceptions it originates as rainfall. Consequently, the measurement and study of rainfall is in actuality the measurement and study of our potential water supply.

Rainfall studies attempt to derive models, both probabilistic and physical, to describe and forecast the rainfall process. Since the quality of every study is immediately related to the quality of the data used, the need for "good quality" rainfall data has been expressed by all hydrologists. By "good quality" is meant accurate, long and uninterrupted series of rainfall measurements at a range of different time intervals (e.g., hourly, daily, monthly, and yearly data) and for a dense raingage network. Missing

values in the series (due, for example, to failure of the recording instruments or to deletion of a station) is a real handicap to the hydrologic data users. The estimation of these missing values is often desirable prior to the use of the data.

For instance, the South Florida Water Management District prepared a magnetic tape with monthly rainfall data for all rainfall stations in south Florida for use in this study (T. MacVicar, SFWMD, personal communication, May, 1982). The data included values for the period of record at each station, ranging from over 100 years (at Key West) to only a few months at several temporary stations. Approximately one month was required to preprocess these data prior to performing routine statistical and time series analyses. The preprocessing included tasks such as manipulations of the magnetic tape, selection of stations with desirable characteristics (e.g., long period of record, proximity to other stations of interest, few missing values) and a major effort at replacement of missing values that did This effort, in fact, was the motivation for this exist. thesis.

Many different kinds of statistical analyses may be performed on a given data set, e.g., determination of elementary statistical parameters, auto- and crosscorrelation analysis, spectral analysis, frequency analysis, fitting time series models. For routine statistics (e.g., calculation of mean, variance and skewness) missing values

are seldom a problem. But for techniques as common as autocorrelation and spectral analysis missing values can cause difficulties. In multivariate analysis missing values result in "wasted information" when only the overlapping period of the series can be used in the analysis, and in inconsistencies (Fiering, 1968, and Chapter 4 of this thesis) when the incomplete series are used.

In general, two approaches to the problem of missing observations exist. The first consists of developing methods of analysis that use only the available data, the second in developing methods of estimation of the missing observations followed by application of classical methods of analysis.

Monthly rainfall totals are usually calculated as the sum of daily recorded values. Thus, if one or more daily observations are missing the monthly total is not reported for that month. An investigation conducted by the Weather Bureau in 1950 (Paulhus and Kohler, 1952), showed that almost one third of the stations for which monthly and yearly totals were not published had only a few (less than five) days missing. Furthermore, for some of these missing days there was apparently no rainfall in the area as concluded by the rainfall observations at nearby stations. Therefore, in many cases estimation of a few missing daily rainfall values can provide a means for the estimation of the monthly totals.

Statisticians have been most concerned with the problem of handling short record multivariate data with missing observations in some or all of the variables, but no explicit and simple solutions have been given, apart from a few special cases in which the missing data follow certain patterns. A review of these methods is given by Afifi and Elashoff (1956). In the time domain, "the analysis of time series, when missing observations occur has not received a great deal of attention" as Marshall (1980, p. 567) comments, and he proposes a method for the estimation of the autocorrelations using only the observed values. Jones (1980) attempts to fit an ARMA model to a stationary time series which has missing observations using Akaike's Markovian representation and Kalman's recursive algorithm. In the frequency domain, spectral analysis with randomly missing observations has been examined by Jones (1962), Parzen (1963), Scheinok (1965), Neave (1970) and Bloomfield (1970).

In hydrology, the problem of missing observations has not been studied much as Salas et al. (1980) state:

The filling-in or extension of a data series is a topic which has not received a great deal of attention either in this book or elsewhere. Because of its importance, the subject is expected to be paid more attention in the future. (Salas et al., 1980, p. 464)

Simple and "practicable" methods for the estimation of missing rainfall values for large scale application were proposed by Paulhus and Kohler (1952), for the completion of the rainfall data published by the Weather Bureau. The

study was initiated after numerous requests of the climatological data users. Beard (1973) adopted a multisite stochastic generation technique to fill-in missing streamflow data, and Kottegoda and Elgy (1977) compared a weighted average scheme and a multivariate method for the estimation of missing data in monthly flow series. Hashino (1977) introduced the "concept of similar storm" for the estimation of missing rainfall sequences. Although the same methods of estimation can be applied to both rainfall and runoff series, a specific method is not expected to perform equally well when applied to the two different series due mainly to the different underlying processes. This is true even for rainfall series from different geographical regions, since their distributions may vary greatly as shown in Fig. 1.1.

This analysis will use monthly rainfall data from four south Florida stations. First, a frequency analysis of the missing observations has been performed and their typical pattern has been identified. In this work the term "missing observations" is used for a sequence of missing monthly values restricted to less than twelve, so that unusual cases of lengthy gaps (a year or more of missing values) is avoided since they do not reflect the general situation.

# Frequency Analysis of Missing Observations in the South Florida Monthly Rainfall Records

An analysis of the monthly rainfall series of 213 stations of the South Florida Water Management District



Fig. 1.1. Monthly distribution of fainfall in the United States (after Linsley R.K., Kohler M.A. and Paulhus J.L., Hydrology for Engineers, 1975, McGraw-Hill, 2nd. edition p. 90)

σ

(SFWMD) gave the results shown on Table 1.1. Figure 1.2 shows the probability density function (pdf) plot of the percent m of missing values, f(m), which is defined as the ratio of the probability of occurrence over an interval to the length of that interval (column 4 of Table 1.1). The shape of the pdf f(m) suggests the fit by an exponential distribution

$$f(m) = \lambda e^{-\lambda m}$$
(1.1)

where  $\lambda$  is the parameter of the distribution calculated as the inverse of the expected value of m, E(m);

$$E(m) = \Sigma p(m_i) m_i \qquad (1.2)$$

where  $p(m_i)$  is the probability of having  $m_i$  percent of missing values. The mean value of the percentage of missing values is  $\overline{m} = E(m) = 13.663$ , and therefore the fitted exponential pdf is

$$f(m) = 0.073 e^{-0.073m}$$
(1.3)

which gives an interesting and unexpectedly good fit as shown by Fig. 1.2 and column 5 of Table 1.1

The question now arises as to whether the missing values within a record follow a certain pattern. In



Fig. 1.2. Probability density function, f(m), of the percentage of missing values. Based on 213 stations,  $\overline{m} = 13.663$ %.

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1	2	3	4	5
% of Missing Values	% of Stations	Cumulative % of Stations	Empirical pdf	Fitted Exponential pdf
0-5	30.52	30.52	0.061	0.061
5-10	21.12	51.64	0.042	0.042
10-15	14.55	66.19	0.029	0.029
15-20	13.61	79.80	0.027	0.020
20-25	6.10	85.90	0.012	0.014
25-30	3.29	89.10	0.007	0.010
30-35	1.88	91.70	0.004	0.007
35-40	0.94	92.01	0.002	0.005
40-45	2.35	94.36	0.005	0.003
45-50	2.82	97.18	0.006	0.002
50 <b>-</b> 55	0.47	97.65	0.001	0.002
55 <b>-</b> 60	0.47	98.12	0.001	0.001
60-65	1.41	99.53	0.003	0.001
65-70	0.47	100.00	0.001	0.001

Table 1.1. Frequency Distribution of the Percent of Missing Values in 213 South Florida Monthly Rainfall Records.

particular, if the occurrence of a gap is viewed as an "event" then the distribution of the interevent times (sizes of the interevents) and of the durations of the events (sizes of the gaps) may be examined.

The probability distribution of the size of the interevents (number of values between two successive gaps) has been studied for four "typical" stations of the SFWMD, as far as length of the record, distribution and percent of missing values is concerned. These four stations are:

MRF 6018, Titusville 2W, 1901-1981, 7.5% missing MRF 6021, Fellsmere 4W, 1911-1979, 9.3% missing MRF 6029, Ocala, 1900-1981, 4.4% missing MRF 6005, Plant City, 1892-1981, 8.6% missing

A derived pdf for the four stations combined and the fitted exponential pdf are shown in Fig. 1.3. The mean size of the interevent,  $\overline{\tau}$ , is 19.03 months; therefore, the fitted exponential distribution is

$$f(\tau) = 0.053 e^{-0.053\tau}$$
(1.4)

Also, the probability distribution of the size of the gaps (number of values missing in each gap) has also been studied for the same four stations. These have been treated as discrete distributions since the size of the gap (k = 1, 2,. . ., 11) is small as compared to the interevent times. A probability distribution for the four stations combined is then derived, which is also the discrete probability mass function (pmf). This plot is shown in Fig. 1.4 and suggests either a Poisson distribution or a discretized exponential.



Fig. 1.3. Probability density function,  $f(\tau)$ , of the interevent size. Based on four stations.

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Fig. 1.4. Probability density, f(k), and mass function, p(k), of the gap size. Based on four stations.

The mean value  $\bar{k}$  is 2.237, which is also the parameter  $\lambda$  of the Poisson distribution. The Poisson distribution

$$f(k) = \frac{e^{-\lambda} \lambda^{k}}{k!}$$
(1.5)

is nonzero at k = 0 and does not fit the peak of the empirical point very well at k = 1 (it gives a value of 0.24 instead of the actual 0.53). The fitted continuous exponential pdf shown in Fig. 1.4 gives a better fit in general but also implies a nonzero probability for a gap size near zero. To overcome this problem and to discretize the continuous exponential pdf, the area (probability) under the exponential curve between zero and 1.5 is assigned to k = 1, ensuring a zero probability at k = 0. Areas (probabilities) assigned to values of k > 1 are centered around those points. The fitted discretized exponential and the Poisson are also shown in Fig. 1.4.

The distributions of the size of the gaps (k) and of the size of interevents ( $\tau$ ) will be used to generate randomly distributed gaps in a complete record. Suppose that we have a complete record and desire to remove randomly m percent missing values. If the mean size of the gap ( $\bar{k}$ ) is assumed constant, the mean size of interevent ( $\bar{\tau}$ ) must vary, decreasing as the percent of missing values increases. Let N denote the total number of values in the record, m the where

$$R^{2} = \rho_{1}\phi_{1} + \rho_{2}\phi_{2} + \dots + \rho_{p}\phi_{p}$$
(3.8)

is called the multiple coefficient of determination and represents the fraction of the variance of the series that has been explained through the regression.

If we denote by  $\phi_{kj}$  the jth coefficient in an autoregressive process of order k, then the last coefficient  $\phi_{kk}$  of the model is called the partial autocorrelation coefficient. Estimates of the partial autocorrelation coefficients  $\phi_{11}$ ,  $\phi_{22}$ ,  $\ldots$ ,  $\phi_{pp}$  may be obtained by fitting to the series autoregressive processes of successively higher order, and solving the corresponding Yule-Walker equations. The partial autocorrelation function  $\phi_{kk}$ , k = 1, 2, . . ., p may also be obtained recursively by means of Durbin's relations (Durbin, 1960)

$$\phi_{k+1,k+1} = [r_{k+1} - \sum_{j=1}^{k} \phi_{k,j} r_{k+1-j}]/[1 - \sum_{j=1}^{k} \phi_{k,j} r_{j}]$$
(3.9)

$$\phi_{k+1,j} = \phi_{k,j} - \phi_{k+1,k+1} \phi_{k,k-j+1}$$
 j = 1, 2, . . , k

It can be shown (Box and Jenkins, 1976, p. 55) that the autocorrelation function of a stationary AR(p) process is a mixture of damped exponential and damped sine waves, infinite in extent. On the other hand, the partial autocorrelation function  $\phi_{kk}$  is nonzero for  $k \leq p$  and zero for k > p. The plot of autocorrelation and partial autocorrelation functions of the series may be used to identify the kind and the order of the model that may have generated it (identification of the model).

# Moving Average Models

In a moving average model the deviation of the current value of the process from the mean is expressed as a finite sum of weighted previous shocks a's. Thus a moving average process of order q can be written as:

$$\tilde{z}_{t} = a_{t} - \theta_{1}a_{t-1} - \theta_{2}a_{t-2} - \dots - \theta_{q}a_{t-q}$$
 (3.10)

or

$$\hat{z}_{t} = \theta (B) a_{t}$$
(3.11)

where

$$\theta(B) = 1 - \theta_1 B - \theta_2 B^2 - \dots - \theta_q B^q$$
 (3.12)

is the moving average operator of order q. An MA(q) model contains (q+2) parameters,  $\mu$ ,  $\theta_1$ ,  $\theta_2$ , ...,  $\theta_q$ ,  $\sigma_a^2$  to be estimated from the data.

From the definition of stationarity (see Appendix A) it follows that an MA(q) process is always stationary since  $\theta(B)$  is finite and thus converges for |B| < 1. But for an MA(q) process to be invertible the q moving average coefficients  $\theta_1$ ,  $\theta_2$ ,  $\ldots$ ,  $\theta_q$  must be chosen so that  $\theta^{-1}(B)$ converges on or within the unit circle, in other words the characteristic equation  $\theta(B) = 0$  must have its roots outside the unit circle.

By multiplying equation (3.10) by  $\tilde{z}_{t-k}$  and taking expected values on both sides we define the autocovariance at lag k:

$$\gamma_{k} = E \left[ \left( a_{t} - \theta_{1} a_{t-1} - \dots - \theta_{q} a_{t-q} \right) \left( a_{t-k} - \theta_{1} a_{t-k-1} - \dots - \theta_{q} a_{t-k-q} \right) \right]$$

$$(3.13)$$

which gives

$$\gamma_{0} = (1 + \theta_{1}^{2} + \theta_{2}^{2} + \dots + \theta_{q}^{2}) \sigma_{a}^{2} \qquad k = 0$$
 (3.14)

$$\gamma_{k} = (-\theta_{k} + \theta_{1}\theta_{k+1} + \theta_{2}\theta_{k+2} + \dots + \theta_{q-k}\theta_{q})\sigma_{a}^{2},$$
  

$$k = 1, 2, \dots, q \qquad (3.15)$$

$$\gamma_k = 0 \qquad k > q \qquad (3.16)$$

By substituting in equation (3.15) the value of  $\sigma_a^2$  from equation (3.14) we obtain a set of q nonlinear equations for  $\theta_1, \theta_2, \ldots, \theta_q$  in terms of  $\rho_1, \rho_2, \ldots, \rho_q$ .

$$\rho_{k} = \frac{-\theta_{k} + \theta_{1}\theta_{k+1} + \theta_{2}\theta_{k+2} + \dots + \theta_{q-k}\theta_{q}}{1 + \theta_{1}^{2} + \dots + \theta_{q}^{2}} , \quad k=1,2,\dots,q$$
(3.17)

These equations are analogous to the Yule-Walker equations for an autoregressive process, but they are not linear and so must be solved iteratively for the estimation of the moving average parameters  $\theta$ , resulting in estimates that may not have high statistical efficiency. Again it was shown by Wold (1938) that these parameters may need corrections (e.g., to fit better the correlogram as a whole and not only the first q correlation coefficients), and that there may exist several, at most 2<sup>q</sup> solutions, for the parameters of the moving average scheme corresponding to an assigned correlogram  $\rho_1$ ,  $\rho_2$ , ...,  $\rho_q$ . However, only those  $\theta$ 's are acceptable which satisfy the invertibility conditions.

From equation (3.14) an estimate for the white noise variance  $\sigma^2_{\rm a}$  may be obtained

$$\sigma_{a}^{2} = \frac{\sigma_{z}^{2}}{1 + \theta_{1}^{2} + \theta_{2}^{2} + \dots + \theta_{q}^{2}}$$
(3.18)

According to the duality principle (see Appendix A) an invertible MA(q) process can be represented as an AR process of infinite order. This implies that the partial autocorrelation function  $\phi_{kk}$  of an MA(q) process is infinite in extent. It can be estimated after tedious algebraic manipulations from the Yule-Walker equations by substituting  $\rho_k$  as functions of  $\theta$ 's for k < q and  $\rho_k = 0$  for k > q. So, in contrast to a stationary AR(p) process, the autocorrelation function of an invertible MA(q) process is finite and cuts off after lag q, and the partial autocorrelation function is infinite in extent, dominated by damped exponentials and damped sine waves (Box and Jenkins, 1976).

# Mixed Autoregressive-Moving Average Models

In practice, to obtain a parsimonious parameterization, it will sometimes be necessary to include both autoregressive and moving average terms in the model. A mixed autoregressive-moving average process of order (p,q), ARMA(p,q), can be written as

$$\tilde{z}_{t} = \phi_{1}\tilde{z}_{t-1} + \dots + \phi_{p}\tilde{z}_{t-p} + a_{t} - \theta_{1}a_{t-1} - \dots - \theta_{q}a_{t-q}$$
(3.19)

or

$$\phi(B) \dot{z}_{t} = \theta(B) a_{t}$$
(3.20)

with (p+q+2) parameters,  $\mu$ ,  $\theta_1$ , ...,  $\theta_q$ ,  $\phi_1$ , ...,  $\phi_p$ ,  $\sigma_a^2$  to be estimated from the data.

An ARMA(p,q) process will be stationary provided that the characteristic equation  $\phi(B) = 0$  has all its roots outside the unit circle. Similarly, the roots of  $\theta(B) = 0$  must lie outside the unit circle for the process to be invertible. By multiplying equation (3.19) by  $\overset{\sim}{z}_{t-k}$  and taking expectations we obtain

$$\gamma_{k} = \phi_{1} \gamma_{k-1} + \dots + \phi_{p} \gamma_{k-p} + \gamma_{za}(k) - \theta_{1} \gamma_{za}(k-1) - \dots$$

$$- \theta_{q} \gamma_{za}(k-q) \qquad (3.21)$$

where  $\gamma_{za}(k)$  is the cross covariance function between z and a, defined by  $\gamma_{za}(k) = E[\hat{z}_{t-k}a_t]$ . Since  $z_{t-k}$  depends only on shocks which have occurred up to time t-k, it follows that

$$\gamma_{za}(k) = 0 \qquad k > 0$$

$$\gamma_{za}(k) \neq 0 \qquad k \leq 0$$
(3.22)

and (3.21) implies

or

$$\rho_{k} = \phi_{1}\rho_{k-1} + \phi_{2}\rho_{k-2} + \dots + \phi_{p}\rho_{k-p} \qquad k \ge q + 1 \quad (3.23)$$

 $\phi(B) \rho_k = 0$   $k \ge q + 1 (3.24)$ 

Thus, for the ARMA(p,q) process the first q autocorrelations  $\rho_1$ ,  $\rho_2$ ,  $\ldots$ ,  $\rho_q$  depend directly on the choice of the q moving average paramaters  $\theta$ , as well as on the p autoregressive parameters  $\phi$  through (3.21). The autocorrelations of higher lags  $\rho_k$ ,  $k \ge q + 1$  are determined through the difference equation (3.24) after providing the p starting values  $\rho_{q-p+1}$ , ...,  $\rho_{q}$ . So, the autocorrelation function of an ARMA(p,q) model is infinite in extent, with the first q-p values  $\rho_1$ , ...,  $\rho_{q-p}$  irregular and the others consisting of damped exponentials and/or damped sine waves (Box and Jenkins, 1976; Salas et al., 1980).

### Autoregressive Integrated Moving Average Models

An ARMA(p,q) process is stationary if the roots of  $\phi(B) = 0$  lie outside the unit circle and "explosive nonstationary" if they lie inside. For example, an explosive nonstationary AR(1) model is  $z_t = 2z_{t-1} + a_t$  (the plot of  $z_t$  vs. t is an exponential growth) in which  $\phi(B) = 1 - 2B$ has its root B = 0.5 inside the unit circle. The special case of homogeneous nonstationarity is when one or more of the roots lie on the unit circle. By introducing a generalized autoregressive operator  $\phi_0(B)$ , which has d of its roots on the unit circle, the general model can be written as

$$\phi_0(B) = \phi(B) (1-B)^d \dot{z}_t = \theta(B) a_t$$
 (3.25)

that is

$$\phi(B) w_{+} = \theta(B) a_{+}$$
 (3.26)

where

$$w_{t} = \nabla^{d} \hat{z}_{t} = \nabla^{d} z_{t}$$
(3.27)

and  $\nabla = 1 - B$  is the difference operator. This model corresponds to assuming that the dth difference of the series can be represented by a stationary, invertible ARMA process. By inverting (3.27)

$$z_t = \nabla^{-d} w_t = s^d w_t$$
(3.28)

where S is the infinite summation operator

$$S = 1 + B + B^{2} + ... = (1-B)^{-1} = \nabla^{-1}$$
 (3.29)

Equation (3.28) implies that the nonstationary process  $z_t$  can be obtained by summing or "integrating" the stationary process  $w_t$ , d times. Therefore, this process is called a simple autoregressive integrated moving average process, ARIMA(p,d,q).

It is also possible to take periodic or seasonal differences at lag's of the series, e.g., the 12th difference of monthly series, introducing the differencing operator  $\nabla_s^D$  with the meaning that seasonal differencing  $\nabla_s$  is applied D times on the series. This periodic ARIMA(P,D,Q)<sub>s</sub> model can be written as

$$\Phi(\mathbf{B}^{\mathbf{S}}) \quad \nabla^{\mathbf{D}}_{\mathbf{S}} \mathbf{z}_{\mathsf{t}} = \Theta(\mathbf{B}^{\mathbf{S}}) \mathbf{a}_{\mathsf{t}}$$
(3.30)

The combination of nonperiodic and periodic models leads to the multiplicative ARIMA(p,d,q) x ARIMA(P,D,Q) model which can be written as

$$\phi(B) \quad \Phi(B^{S}) \quad \nabla^{d} \quad \nabla^{D}_{s} \quad z_{t} = \theta(B) \quad \Theta(B^{S}) \quad a_{t}$$
(3.31)

After the model has been fitted to the differenced series an integration should be performed to retrieve the original process. But such an integrated series would lack a mean value since a constant of integration has been lost through the differencing. This is the reason that the ARIMA models cannot be used for synthetic generation of time series, although they are useful in forecasting the deviations of a process (Box and Jenkins, 1976; Salas et al., 1980).

# Transformation of the Original Series

# Transformation to Normality

Most probability theory and statistical techniques have been developed for normally distributed variables. Hydrologic variables are usually assymetrically distributed or bounded by zero (positive variables), and so a transformation to normality is often applied before modeling. Another approach would be to model the original skewed series and then find the probability distribution of the uncorrelated residuals. Care must then be taken to assess the errors of applying methods developed for normal variables to skewed variables, especially when the series are highly skewed, e.q., hourly or daily series. On the other hand, when transforming the original series into normal, biases in the mean and standard deviation of the generated series may occur. In other words, the statistical properties of the transformed series may be reproduced in the generated but not in the original series. An alternative for avoiding biases in the moments of the generated series would be to estimate the moments of the transformed series through the derived relationships between the moments of the skewed and normal series. Matalas (1967) and Fiering and Jackson (1971) describe how to estimate the first two moments of the logtransformed series so as to reproduce the ones of the original series. Mejia et al. (1974) present another approach in order to preserve the correlation structure of the original series.

However, the most widely used approach is to transform the original skewed series to normal and then model the normal series. Several transformations may be applied to the original series, and the transformed series then tested for normality, e.g. the graph of their cumulative distribution should appear as a straight line when it is plotted on normal probability paper. The transformation will be finally chosen that gives the best approximation to normality, e.g., the best fit to a straight line.

Another advantage of transforming the series to normal is that the maximum likelihood estimates of the model
parameters are essentially the same as the least squares estimates, provided that the residuals are normally distributed (Box and Jenkins, 1976, Ch. 7). This facilitates the calculation of the final estimates since they are those values that minimize the sum of squares of the residuals.

Box and Cox (1964) showed how a maximum likelihood and a parallel Bayesian analysis can be applied to any type of transformation family to obtain the "best" choice of transformation from that family. They illustrated those methods for the popular power families in which the observation x is replaced by y, where

$$y = \begin{cases} \frac{x^{\lambda} - 1}{\lambda} , & \lambda \neq 0 \\ \log x , & \lambda = 0 \end{cases}$$
(3.32)

The fundamental assumption was that for some  $\lambda$  the transformed observations y can be treated as independently normally distributed with constant variance  $\sigma^2$  and with expectations defined by a linear model

 $E[y] = A L \tag{3.33}$ 

where A is a known constant matrix and L is a vector of unknown parameters associated with the transformed observations (Box and Cox, 1964).

This transformation has the advantage over the simple power transformation proposed by Tukey (1957)

$$y = \begin{cases} x^{\lambda}, & \lambda \neq 0 \\ \log x, & \lambda = 0 \end{cases}$$
(3.34)

of being continuous at  $\lambda=0$ . Otherwise the two transformations are identical provided, as has been shown by Schlesselman (1971), that the linear model of (3.33) contains a constant term.

Further, Draper and Cox (1969), showed that the value of  $\lambda$  obtained from this family of transformations can be useful even in cases where no power transformation can produce normality exactly. Also, John and Draper (1980) suggested an alternative one-parameter family of transformations when the power transformation fails to produce satisfactory distributional properties as in the case of a symmetric distribution with long tails.

The selection of the exact transformation to normality (zero skewness) is not an easy task, and over-transformation, i.e., transformation of the original data with a large positive (negative) skewness to data with a small negative (positive) skewness, or under-transformation, i.e., transformation of the original data with a large positive (negative) skewness to data with a small positive (negative) skewness, may result in unsatisfactory modeling of the series or in forecasts that are in error. This was the case for the data used by Chatfield and Prothero (1973a), who applied the Box-Jenkins forecasting approach and were dissatisfied with the results, concluding that the Box-Jenkins forecasting procedure is less efficient than other forecasting

methods. They applied a log transform to the data which evidently over-transformed the data, as shown by Box and Jenkins (1973) who finally suggested the approximate transformation  $y = x^{0.25}$ , even though the complicated but precise Box-Cox procedure gave an estimate of  $\lambda = 0.37$  [Wilson (1973)].

Thus, the selection of the normality transformation greatly affects the forecasts, as Chatfield and Prothero (1973b) experienced with their data. They concluded that

. . We have seen that a "small" change in  $\lambda$ from 0 to 0.25 has a substantial effect on the resulting forecasts from model A [ARIMA(1,1,1) x ARIMA(1,1,1)<sub>12</sub>] even though the goodness of fit does not seem to be much affected. This reminds us that a model which fits well does not necessarily forecast well. Since small changes in  $\lambda$  close to zero produce marked changes in forecasts, it is obviously advisable to avoid "low" values of  $\lambda$ , since a procedure which depends critically on distinguishing between fourth-root and logarithmic transformation is fraught with peril. On the other hand a "large" change in  $\lambda$  from 0.25 to 1 appears to have relatively little effect on forecasts. So we conjecture that Box-Jenkins forecasts are robust to changes in the transformation parameter away from zero. . . .[Chatfield and Prothero (1973b) p. 347]

### Stationarity

Most time series occurring in practice exhibit nonstationarity in the form of trends or periodicities. The physical knowledge of the phenomenon being studied and a visual inspection of the plot of the original data may give the first insight into the problem. Usually the length of the series is not long enough, and the detection of

trends or cycles only through the plot of the series is ambiquous. Useful tools for the detection of periodicities are the autocorrelation function and the spectral density function of the series (which is the Fourier transform of the autocorrelation function). If a seasonal pattern is present in the series then the correlogram (plot of the autocorrelation function) will exhibit a sinusoidal appearance and the periodogram (plot of the spectral density function) will show peaks. The period of the sinusoidal function of the correlogram, or the frequency where the peaks occur in the periodogram, can determine the periodic component exactly (Jenkins and Watts, 1968). Another device for the detection of trends and periodicities is to fit some definite mathematical function, such as exponentials, Fourier series or polynomials to the series and then model the residual series, which is assumed to be stationary. More details on the treatment of nonstationary data as well as on the interpretation of the correlogram and periodogram of a time series can be found in textbooks such as Bendat and Piersol (1958), Jenkins and Watts (1968), Wastler (1969), Yevjevich (1972), and Chatfield (1980).

Apart from the approach of removing the nonstationarity of the original series and modeling the residual series with a stationary ARMA(p,q) model, the original nonstationary series can be modeled directly with a simple or seasonally integrated ARIMA model. Actually, the second approach can be viewed as an extension of the first one,

e.g., the nonstationarity is removed through the simple  $(\nabla)$  or seasonal  $(\nabla_s)$  differencing. However, the integrated model cannot be used for generation of data, as has already been discussed.

For many hydrologic applications, one is satisfied with second order or weak stationarity, e.g., stationarity in the mean and variance. Furthermore, weak stationarity and the assumption of normality imply strict stationarity (see Appendix A).

#### Monthly Rainfall Series

## Normalization and Stationarization

Stidd (1953, 1968) suggested that rainfall data have a cube root normal distribution because they are product functions of three variables: vertical motion in the atmosphere, moisture, and duration time. Synthetic rainfall data generated using processes analogous to those operating in nature showed that the exponent required to normalize the distribution is between 0.5 (square root) and 0.33 (cubic root) for different types of rainfall (Stidd, 1970).

The square root transformation has been extensively used for the approximate normalization of monthly rainfall series (see Table Cl2 of Appendix C) with satisfactory results: Delleur and Kavvas (1978), Salas <u>et al</u>. (1980), Ch. 5, Roesner and Yevjevich (1966). However, Hinkley (1977) used the exact Box-Cox transformation for monthly rainfall series. Although, Asley <u>et al</u>. (1977) have developed an efficient algorithm for the estimation of  $\lambda$  along with other parameters in an ARIMA model, it seems that the exact value of  $\lambda$  is not more reliable than the approximate one  $\lambda = 0.5$ (Chatfield and Prothero, 1973b). The reasons for this follow.

First, Chatfield and Prothero (1973b) used the Box-Cox procedure to evaluate the exact transformation of their data. They obtained estimates  $\hat{\lambda} = 0.24$  using all the data (77 observations),  $\hat{\lambda} = 0.34$  using the first 60 observations and  $\hat{\lambda} = 0.16$  excluding the first year's data. Therefore, it is logical to infer that even if the complicated Box-Cox procedure for the incomplete rainfall record is used, the missing values may be enough to give a spurious  $\lambda$ , which is not "more exact" than the value of 0.5 used in practice.

Second, we may also notice that the use of either  $\lambda = 0.33$  (cubic root) or  $\lambda = 0.5$  (square root) is not expected to greatly affect the forecasts since, according to Chatfield and Prothero (1973b), the Box-Jenkins forecasts are not too sensitive to changes of  $\lambda$  for  $\lambda > 0.25$ .

Monthly rainfall series are nonstationary. The variation in the mean is obvious since generally the expected monthly rainfall value for January is not the same as that of July. Although the variation of the standard deviation is not so easy to visualize, calculations show that months with higher mean usually have higher standard deviation. Thus, each month has its own probability

distribution and its own statistical parameters resulting in monthly series that are nonstationary.

By introducing the concept of circular stationarity as developed by Hannan (1960) and others (see Appendix A for definition), the periodic monthly rainfall series can be considered <u>not</u> as nonstationary but circular stationary, since circular stationarity suggests that the probability distribution of rainfall in a particular month is the same for the different years. Then, the monthly rainfall series is composed of a circularly stationary (periodic) component and a stationary random component.

The time-series models currently used in hydrology are fitted to the stationary random component, so the circularly stationary component must be removed before modeling. This last component appears as a sinusoidal component in the autocorrelation function (with a 12-month period) or as a discrete spectral component in the spectrum (peak at the frequency 1/12 cycle per month). Usually several subharmonics of the fundamental 12-month period are needed to describe all the irregularities present in the autocorrelation function and spectral density function, since in nature the periodicity does not follow an ideal cosine The use of a Fourier function with a 12-month period. series approach for the approximation of the periodic component of monthly rainfall and monthly runoff series has been illustrated by Roesner and Yevjevich (1966).

Kavvas and Delleur (1975) investigated three methods of removal of periodicities in the monthly rainfall series: nonseasonal (first-lag) differencing, seasonal differencing (12-month difference), and removal of monthly means. They worked both analytically and empirically using the rescaled (divided by the monthly standard deviation) monthly rainfall square roots for fifteen Indiana watersheds. They concluded that "all the above transformations yield hydrologic series which satisfy the classical second-order weak stationarity conditions. Both seasonal and nonseasonal differencing reduce the periodicity in the covariance function but distort the original spectrum, thus making it impractical or impossible to fit an ARMA model for generation of synthetic monthly series. The subtraction of monthly means removes the periodicity in the covariance and the amount of nonstationarity introduced is negligible for practical purposes." (Kavvas and Delleur, 1975, p. 349.) In other words, they concluded that the best way for modeling monthly rainfall series is to remove the seasonality (by subtracting the monthly means and dividing by the standard deviations of the normalized series) and then use a stationary ARMA(p,q) model to model the stationary normal residuals.

#### Modeling of Normalized Series

It is assumed that the nonstationarities due to longterm trends are removed before any operation. Then the appropriate transformation is applied to the data in

order to obtain an approximately normal distribution. For monthly rainfall series experience has shown that the best practical transformation is the square root transformation, as has already been discussed. What remains is the modeling of the normalized series with one of the following models: stationary ARMA(p,q), simple nonstationary ARIMA(p,d,q), seasonal nonstationary ARIMA(P,D,Q)<sub>s</sub>, or multiplicative ARIMA(p,d,q)x(P,D,Q)<sub>s</sub> model.

Delleur and Kavvas (1978) fitted different models to the monthly rainfall series of 15 basins in Indiana and compared the results. They studied the models: ARIMA (0,0,0), ARIMA(1,0,1), ARIMA(1,1,1), ARIMA(1,1,1)<sub>12</sub>, and ARIMA(1,0,0)x(1,1,1)<sub>12</sub> on the square-root trans-They concluded that from the nonseasonal formed series. ARIMA models, ARMA(1,1) "emerged as the most suitable for the generation and forecasting of monthly rainfall series." The goodness-of-fit tests applied on the residuals were the portemanteau lack of fit test (see Appendix A) of Box and Pierce (1970) and the cumulative periodogram test (Box and Jenkins, 1976, p. 294). The ARMA(1,1) model passed both tests in all cases studied. From the nonseasonal models,  $ARIMA(1,0,0)x(1,1,1)_{12}$  also passed the goodness-of-fit tests in all cases, but they stress that this model "has only limited use in the forecasting of monthly rainfall series since it does not preserve the monthly standard deviations." As far as forecasts are concerned, they showed that "the forecasts by the several models follow each other very

closely and the forecasts rapidly tend to the mean of the observed rainfall square roots (which is the forecast of the white noise model)."

## CHAPTER 4

## MULTIVARIATE STOCHASTIC MODELS

#### Introduction

For univariate stochastic models the sequence of observations under study is assumed independent of other sequences of observations and so is studied by itself (single or univariate time series). However, in practice there is always an interdependence among such sequences of observations, and their simultaneous study leads to the concept of multivariate statistical analysis. For example, a rainfall series of one station may be better modeled if its correlation with concurrent rainfall series at other nearby stations is incorporated into the model. Multiple time series can be divided into two groups: (1) multiple time series at several points (e.g., rainfall series at different stations, streamflow series at various points of a river), and (2) multiple series of different kinds at one point (e.g., rainfall and runoff series at the same station). In general, both kinds of multiple time series are studied simultaneously, and their correlation and cross-correlation structure is used for the construction of a model that better describes all these series. The parameters of this so called multivariate stochastic model are calculated such

that the correlation and cross-correlation structure of the multiple measured series are preserved in the multiple series generated by the model.

The multivariate models that will be presented in this chapter have been developed and extensively used for the generation of synthetic series. How these models can be adapted and used for filling in missing values will be discussed in chapter 5.

#### General Multivariate Regression Model

The general form of a multivariate regression model is

$$Y = A X + B H \tag{4.1}$$

where Y is the vector of dependent variables, X the vector of independent variables, A and B matrices of regression coefficients, and H a vector of random components. The vectors Y and X may consist of either the same variable at different points (or at different times) or different variables at the same or different points (or at different times).

For convenience and without loss of generality all the variables are assumed second order stationary and normally distributed with zero mean and unit variance. Transformations to accomplish normality have been discussed in Chapter 3. A random component is superimposed on the model to account for the nondeterministic fluctuations.

In the above model, the dependent and independent variables must be selected carefully so that the most

information is extracted from the existing data. A good summary of the methods for the selection of independent variables for use in the model is given in Draper and Smith (1966). Most popular is the stepwise regression procedure in which the independent variables are ranked as a function of their partial correlation coefficients with the dependent variable and are added to the model, in that order, if they pass a sequential F test.

The parameter matrices A and B are calculated from the existing data in such a way that important statistical characteristics of the historical series are preserved in the generated series. This estimation procedure becomes cumbersome when too many dependent and independent variables are involved in the model, and several simplifications are often made in practice. On the other hand, restrictions have to be imposed on the form of the data, as we shall see later, to ensure the existence of real solutions for the matrices A and B.

#### Multivariate Lag-One Autoregressive Model

If only one variable (e.g., rainfall at different stations) is used in the analysis then the model of equation (4.1) becomes a multivariate autoregressive model. Since in the rest of this chapter we will be dealing only with one variable (rainfall) which has been transformed to normal and second order stationary, the vectors Y and X are replaced by the vector Z for a notation consistent with the

univariate models. Matalas (1967) suggested the multivariate lag-one autoregressive model

$$Z_{t} = A Z_{t-1} + B H_{t}$$
(4.3)

where  $Z_t$  is an (mxl) vector whose ith element  $z_{it}$  is the observed rainfall value at station i and at time t, and the other variables have been described previously.

Such a model can be used for the simultaneous generation of rainfall series at m different stations. The correlation and cross-correlation of the series is incorporated in the model through the parameters A and B.

The matrices A and B are estimated from the historical series so that the means, standard deviations and autocorrelation coefficients of lag-one for all the series, as well as the cross-correlations of lag-zero and lag-one between pairs of series are maintained.

Let  ${\rm M}_{\mbox{\scriptsize 0}}$  denote the lag-zero correlation matrix which is defined as

$$M_0 = E[Z_t \ Z_t^T]$$
(4.4)

Then a diagonal element of  $M_0$  is  $E[z_{i,t} z_{i,t}] = \rho_{ii}(0) = 1$ (since  $Z_t$  is standardized) and an off diagonal element (i,j) is  $E[z_{i,t} z_{j,t}] = \rho_{ij}(0)$  which is the lag-zero cross correlation between series  $\{z_i\}$  and  $\{z_j\}$ . The matrix  $M_0$  is symmetric since  $\rho_{ij}(0) = \rho_{ji}(0)$  for every i, j. Let  $M_1$  denote the lag-one correlation matrix defined as

$$M_{1} = E[Z_{t} \ Z_{t-1}^{T}]$$
(4.5)

A diagonal element of  $M_1$  is  $E[z_{i,t} z_{i,t-1}] = \rho_{ii}(1)$  which is the lag-one serial correlation coefficient of the series  $\{z_i\}$ , and an off-diagonal element (i,j) is  $E(z_{i,t} z_{j,t-1}) = \rho_{ij}(1)$  which is the lag-one cross-correlation between the  $\{z_i\}$  and  $\{z_j\}$  series, the latter lagged behind the former. Since in general  $\rho_{ij}(1) \neq \rho_{ji}(1)$  for  $i \neq j$  the matrix  $M_1$  is not symmetric.

After some algebraic manipulations (see Appendix B) the coefficient matrices A and B are obtained as solutions to the equations

$$A = M_1 M_0^{-1}$$
 (4.6)

$$B B^{T} = M_{0} - M_{1} M_{0}^{-1} M_{1}^{T}$$
(4.7)

where  $M_0^{-1}$  is the inverse of  $M_0$ , and  $M_1^T$  the transpose of  $M_1$ . The correlation matrices  $M_0$  and  $M_1$  are calculated from the data. Then an estimate of the matrix A is given directly by equation (4.6), and an estimate for B is found by solving equation (4.7) by using a technique of principal component analysis (Fiering, 1964) or upper triangularization (Young, 1968). For more details on the solution of equation (4.7) see Appendix B.

## Comments on Multivariate AR(1) Model

## Assumption of Normality and Stationarity

We have assumed that all random variables involved in the model are normal. The assumption of a multivariate normal distribution is convenient but not necessary. It has been shown (Valencia and Schaake, 1973) that the multivariate AR(1) model preserves first and second order statistics regardless of the underlying probability distributions.

Several studies have been done using directly the original skewed series. Matalas (1967) worked with lognormal series and constructed the generation model so that it preserves the historical statistics of the log-normal process. Mejia et al. (1974) showed a procedure for multivariate generation of mixtures of normal and log-normal variables. Moran (1970) indicated how a multivariate gamma process may be applied, and Kahan (1974) presented a method for the preservation of skewness in a linear bivariate regression model. But in general, the normalization of the series prior to modeling is more convenient, especially when the series have different underlying probability distribu-In such cases different transformations are applied tions. on the series, and that combination of transformations is kept which yields minimum average skewness. Average skewness is the sum of the skewness of each series divided by the number of series or number of stations used. This operation is called finding the MST (Minimum Skewness

Transformation) and results in an approximately multivariate normal distribution (Young and Pisano, 1968).

We have also assumed that all variables are standardized, e.g., have zero mean and unit variance. This assumption is made without loss of generality since the linear transformations are preserved through the model. On the other hand this transformation becomes necessary when modeling periodic series since by subtracting the periodic means and dividing by the standard deviations we remove almost all of the periodicity.

If the data are not standardized,  $M_0$  and  $M_1$  represent the lag-zero and lag-one covariance matrices (instead of correlation matrices), respectively. If S denotes the diagonal matrix of the standard deviations and  $R_0$ ,  $R_1$  the lag-zero and lag-one correlation matrices then

$$M_0 = S R_0 S \tag{4.8}$$

and

 $M_{1} = S R_{1} S \tag{4.9}$ 

When we standardize the data the matrix S is an identity matrix and  $M_0$ ,  $M_1$  become the correlation matrices  $R_0$  and  $R_1$ respectively. Thus, one other advantage of standardization is that we work with correlation matrices whose elements are less than unity and the computations are likely to be more stable (Pegram and James, 1972).

# Cross-Correlation Matrix M1

Notice that the lag-one correlation matrix  $M_1$  has been defined as  $M_1 = E[Z_t Z_{t-1}^T]$  which contains the lag-one cross-correlations between pairs of series but having the second series lagged behind the first one. Following this definition the lag-minus-one correlation matrix will be

$$M_{-1} = E[Z_{t-1} \ Z_{t}^{T}]$$
(4.10)

and it will contain the lag-one correlations having now the second series lagged ahead of the first one. It is easy to show that  $M_{-1}$  is actually the transpose of  $M_1$ :

$$M_{-1} = E[Z_{t-1} \ Z_{t}^{T}] = E[(Z_{t} \ Z_{t-1}^{T})^{T}] = M_{1}^{T}$$
(4.11)

Care then must be taken so that there is a consistency between the equation used to calculate matrix A and the way that the cross-correlation coefficients have been calculated. Such an inconsistency was present in the numerical multisite package (NMP) developed by Young and Pisano (1968) and was first corrected by O'Connell (1973) and completely corrected and improved by Finzi et al. (1974, 1975).

#### Incomplete Data Sets

In practice, hydrologic series at different stations are unlikely to be concurrent and of equal length. With lag-zero auto- and cross-correlation coefficients calculated from the incomplete data sets, the lag-zero correlation matrix  $M_0$  obtained may not be positive semidefinite, and, its inverse  $M_0^{-1}$  needed for the calculation of matrix A thus may have elements that are complex numbers. Also, a necessary and sufficient condition for a real solution of matrix B is that  $C = M_0 - M_1 M_0^{-1} M_1^T$  is a positive semi-definite matrix (see Appendix B).

When all of the series are concurrent and complete then  $M_0$  and C are both semidefinite matrices [Valencia and Schaake, 1973], and the generated synthetic series are real numbers. When the series are incomplete there is no guarantee that real solutions for the matrices A and B exist causing the model of Matalas (1967) to be conditional on  $M_0$ and C being positive semidefinite [Slack, 1973].

Several techniques have been proposed which use the incomplete data sets but guarantee the posite semidefiniteness of the correlation matrices. Fiering (1968) suggested a technique that can be used to produce a positive semi-definite correlation matrix  $M_0$ . If  $M_0$  is not positive semidefinite then negative eigenvalues may occur and hence negative variables, since the eigenvalues are variances in the principal component system. In this technique, the eigenvalues of the original correlation matrix are calculated. If negative eigenvalues are encountered, an adjustment procedure is used to eliminate them (thereby altering the correlation matrix,  $M_0$  [Fiering, 1968]).

A correlation matrix is called consistent if all its eigenvalues are positive. But consistent estimates of the correlation matrices  $M_0$  and  $M_1$  do not guarantee that C will also be consistent.

Crosby and Maddock (1970) proposed a technique that is suitable only for monotone data (data continuous in collection to the present but having different starting times). This technique produces a consistent estimate of the matrix  $M_0$  as well as of the matrix C, and is based on the maximum likelihood technique developed by Anderson (1957).

Valencia and Schaake (1973) developed another technique. They estimate matrices A and B from the equations

$$A = M_{1} M_{01}^{-1}$$
(4.12)

$$B B^{T} = M_{02} - M_{1} M_{01}^{-1} M_{1}^{T}$$
(4.13)

where  $M_{01}$  is the lag-zero correlation matrix  $M_0$  computed from the first (N-1) vectors of the data, and  $M_{02}$  is computed from the last (N-1) vectors, where N is the number of data points (number of times sampled) in each of the n series.

#### Further Simplification

Sometimes in practice, the preservation of the lagzero and lag-one autocorrelations and the lag-zero cross-correlations is enough. In such cases, i.e., when the lag-one cross-correlations are of no interest, a nice simplification can be made due to Matalas (1967, 1974). He defined matrix A as a diagonal matrix whose diagonal elements are the lag-one auto-correlation coefficients. With A defined as above, the lag-one cross-correlation of the generated series ( $\rho'_{ij}(1)$ ) can be shown to be the product of the lag-zero cross-correlation ( $\rho_{ij}(0)$ ) and the lag-one auto-correlation of the series different than the actual lag-one cross-correlation ( $\rho_{ij}(1)$ ).

$$\rho_{ij}(l) = \rho_{ij}(0) \rho_{ii}(l)$$
 (4.14)

By using  $\rho_{ij}(1)$  of equation (4.14) in place of the actual  $\rho_{ij}(1)$ , thus avoiding the actual computation of  $\rho_{ij}(1)$  from the data, the desired statistical properties of the series are still preserved.

#### Higher Order Multivariate Models

The order p of a multivariate autoregressive model could be estimated from the plots of the autocorrelation and partial autocorrelation functions of the series (Salas <u>et al.</u>, 1980) as an extension of the univariate model identification, which is already a difficult and ambiguous task. However, in practice first and second order models are usually adequate and higher order models should be avoided (Box and Jenkins, 1976). In any case, the multivariate multilag autoregressive model of order p takes the form

$$Z_{t} = \sum_{k=1}^{p} A_{k} Z_{t-k} + B H_{t}$$
(4.15)

and the matrices  $A_1, A_2, \dots A_p$ , B are the solutions of the equations

$$M_{i} = \sum_{k=1}^{p} A_{k} M_{i-k} , i = 1, 2, ..., p \qquad (4.16)$$

$$B B^{T} = M_{0} - \sum_{k=1}^{p} A_{k} M_{k}^{T}$$

$$(4.17)$$

where  $M_{l}$  is the lag-l correlation matrix. Equation (4.16) is a set of p matrix equations to be solved for the matrices  $A_{1}, A_{2}, \ldots, A_{p}$ , and matrix B is obtained from (4.17) using techniques already discussed. Here, the assumption of diagonal A matrices becomes even more attractive. For a multivariate second-order AR process the above simplification is illustrated in Salas and Pegram (1977) where the case of periodic (not constant) matrix parameters is also considered.

O'Connell (1974) studied the multivariate ARMA(1,1) model

$$Z_{t} = A Z_{t-1} + B H_{t} - C H_{t-1}$$
 (4.18)

where A, B, and C are coefficient matrices to be determined

from the data. Specifically they are solutions of the system of matrix equations

$$B B^{T} + C C^{T} = S$$

$$C B^{T} = T$$
(4.19)

where S and T are functions of the correlation matrices  $M_0$ ,  $M_1$  and  $M_2$ . Methods for solving this system are proposed by O'Connell (1974).

Explicit solutions for higher order multivariate ARMA models are not available and Salas <u>et al</u>. (1980) propose an approximate multivariate ARMA(p,q) model.

## CHAPTER 5

## ESTIMATION OF MISSING MONTHLY RAINFALL VALUES--A CASE STUDY

#### Introduction

This section compares and evaluates different methods for the estimation of missing values in hydrological time series. A case study is presented in which four of the simplified methods presented in Chapter 2 have been applied to a set of four concurrent 55 year monthly rainfall series from south Florida and the results compared. Also a recursive method for the estimation of missing values by the use of a univariate or multivariate stochastic model has been proposed and demonstrated. The theory already presented in Chapters 2, 3 and 4 is supplemented whenever needed.

#### Set Up of the Problem

The monthly rainfall series of four stations in the South Florida Water Management District (SFWMD) have been used in the analysis. These stations are:

Station A : MRF6038, Moore Haven Lock 1
Station 1 : MRF6013, Avon Park
Station 2 : MRF6093, Fort Myers WSO Ap.
Station 3 : MRF6042, Canal point USDA.

For convenience the four stations will sometimes be addressed as A, 1, 2, 3 instead of their SFWMD identification numbers 6038, 6013, 6093 and 6042, respectively. Their locations are shown in the map of Fig. 5.1. Station A in the center is considered as the interpolation station (whose missing values are to be estimated) and the other three stations 1, 2 and 3 as the index stations. Care has been taken so that the three index stations are as close and as evenly distributed around the interpolation station as possible.

This particular set of four stations was selected because it exhibits many desired and convenient properties:

- the stations have an overlapping period of 55 years
   (1927-1981),
- (2) for this 55 year period the record of the interpolation station (station A) is complete (no missing values),
- (3) the three index stations have a small percent of missing values for the overlapping period (station 1: 2.7% missing, station 2: complete, and station 3: 1.2% missing values).

The 55 year length of the records is considered long enough to establish the historical statistics (e.g., monthly mean, standard deviation and skewness) and provides a monthly series of a satisfactory length (660 values) for fitting a univariate or multivariate ARMA model.



The four south Florida rainfall stations Fig. 5.1. used in the analysis.

- A: 6038, Moore Haven Lock 1
- 1: 6013, Avon Park
- 2: 6093, Fort Myers WSO AP. 3: 6042, Canal Point USDA

The completeness of the series of the interpolation station permits the random generation of gaps in the series, corresponding to different percentages of missing values, with the method described in Chapter 1. After the missing values have been estimated by the applied models, the gaps are in-filled with the estimated values and the statistics of the new (estimated) series are compared with the statistics of the incomplete series and the statistics of the historical (actual) series. Also the statistical closeness of the in-filled (estimated) values to the hidden (actual) values provides a means for the evaluation and comparison of the methods.

When, for the estimation of a missing value of the interpolation station, the corresponding value of one or more index stations is also missing the latter is eliminated from the analysis, e.g., only the remaining one or two index stations are used for the estimation. Frequent occurrence of such concurrent gaps in both the interpolation and the index stations would alter the results of the applied method in a way that cannot be easily evaluated (e.g., another parameter such as the probability of having concurrent gaps should be included in the analysis). A small number of missing values in the selected index stations eliminates the possibility of such simultaneous gaps, and thus the effectiveness of the applied estimation procedures can be judged more efficiently.

The statistical properties (e.g., monthly mean, standard deviation, skewness and coefficient of variation) of the truncated (to the 1927-1981 period) original monthly rainfall series for the four stations are shown on Tables C.1, C.2, C.3 and C.4 of Appendix C. Figure 5.2 shows the plot of the monthly means and standard deviations for station A. From these plots we observe that: (1) the plot of monthly means is in agreement with the typical plot for Florida shown in Fig. 1.1, and (2) months with a high mean usually have a high standard deviation. The only exception seems to be the month of January which in spite of its low mean exhibits a high standard deviation and therefore a very high coefficient of variation and an unusually high skewness. A closer look at the January rainfall values of station A shows that the unusual properties for that month are due to an extreme value of 21.4 inches of rainfall for January 1979, the other values being between 0.05 and 6.04 inches.

The three index stations 1, 2 and 3 are at distances 59 miles, 51 miles and 29 miles respectively from the interpolation station A.

#### Simplified Estimation Techniques

#### Techniques Utilized

From the simplified techniques presented in Chapter 2, the following four are applied for the estimation of missing





### monthly rainfall values:

(1) the mean value method (MV)

- (2) the reciprocal distances method (RD)
- (3) the normal ratio method (NR), and
- (4) the modified weighted average method (MWA).

These methods are all deterministic and are applied directly on the available data permitting thus a uniform and objective comparison of the results. The mean value plus random component method has not been included in this thesis.

The above four methods will be applied for five different percentages of missing values: 2%, 5%, 10%, 15% and 20%. These percentages cover almost 80% of all cases encountered in practice as has been shown in Table 1.1 (e.g., 80% of the stations have below 20% missing values). From the same table it can also be seen that almost 30% of the stations have below 5% missing values. Therefore, it would be of interest and practical use if we could generalize the results for the region of below 5% missing values since a large fraction of the cases in practice fall in this region.

The application of the first three methods (MV, RD, NR methods) is straightforward and no further comments need be made. However, some comments on the least squares (LS) method and the modified weighted average (MWA) method are necessary.

## Least Squares Method (LS)

The least squares method although simple in principle involves an enormous amount of calculations, and for that reason it has been excluded from this study. For example, consider the case in which the interpolation station A is regressed on the three index stations 1, 2 and 3. The estimated values will be given by:

$$y' = a + b_1 x_1 + b_2 x_2 + b_3 x_3 + \varepsilon$$
 (5.1)

where a, b<sub>1</sub>, b<sub>2</sub>, b<sub>3</sub> are the regression coefficients calculated from the available concurrent values of all the four variables. There are 12 such regression equations, one for each month. But if it happens that an index station (say, station 3) has a missing value simultaneously with the interpolation station, a new set of 12 regression equations is needed for the estimation, e.g.,

$$y' = a' + b'_1 x_1 + b'_2 x_2 + \varepsilon$$
 (5.2)

Unless this coincidence of simultaneously missing values is investigated manually so that only the needed least squares regressions are performed (Buck, 1960), all the possible combinations of regressions must otherwise be performed. This involves regressions among all the four variables (y;  $x_1$ ,  $x_2$ ,  $x_3$ ), among the three of them (y;  $x_1$ ,  $x_2$ ), (y;  $x_1$ ,  $x_3$ ), (y;  $x_2$ ,  $x_3$ ) and between pairs of them (y;  $x_1$ ), (y;  $x_2$ ), (y;  $x_3$ ), giving overall 7 sets of 12 regression equations. Because the regression coefficients are different for each percentage of missing values (since their calculation is based only on the existing concurrent values) the 84 (7 x 12) regressions must be repeated for each level of missing values (420 regressions overall for this study).

It could be argued that the same 12 regression equations (y;  $x_1$ ,  $x_2$ ,  $x_3$ ) could be kept and a missing values  $x_i$  replaced by its mean  $\overline{x}_i$  or by another estimate  $x'_i$ . In that case equation 5.1 would become

$$y' = a + b_1 x_1 + b_2 x_2 + b_3 x_3' + \varepsilon,$$
 (5.3)

the coefficients of regression a,  $b_1$ ,  $b_2$ ,  $b_3$  remaining unchanged. This in fact can be done, but then the method tested will not be the "pure" least squares method since the results will depend on the secondary method used for the estimation of the missing  $x_i$  values.

The coefficients a,  $b_1$ ,  $b_2$  and  $b_3$  (equation 5.1) of the regression of the {y} series (of station A with 2% missing values) on the series  $\{x_1\}$ ,  $\{x_2\}$  and  $\{x_3\}$  (of stations 1, 2 and 3 respectively) are shown in Table 5.1. In the same table the values of the squared multiple regression coefficient  $R^2$  and the standard deviation of the {y} series are also shown. The numbers in parenthesis show the significance level  $\alpha$  at which the parameters are significant (the percent probability of being nonzero is  $(1-\alpha)$ )100. For

also given. R<sup>2</sup>  $b_1$ а b<sub>2</sub> bz S inches inches 0.0059 0.4994 0.3377 0.8046 0.1271 3.076 JAN (0.9692)(0.2790)(0.0005)(0.0017)(0.0001)0.1355 0.2624 0.0086 0.5345 0.7033 FEB 1.365 (0.5260)(0.0025)(0.9431)(0.0001)(0.0001)0.0052 0.1617 0.3457 0.4507 0.9142 MAR 2.464 (0.0001)(0.9793)(0.0138)(0.0001)(0.0001)0.7388 0.2405 0.2813 0.1919 0.4936 APR 1.818 (0.0273)(0.0156)(0.1132)(0.0001)(0.0458)2.1302 0.4046 -0.0591 0.2186 0.2752 MAY 2.583 (0.0070)(0.0115)(0.7180)(0.1308)(0.0016)1.8765 0.2192 0.1108 0.3339 0.3351 JUN 3.812 (0.4034)(0.0133)(0.0002)(0.1505)(0.1576)2.8601 -0.0345 0.3993 0.1885 0.2005 JUL 3.399 (0.0750)(0.7883)(0.0131)(0.1780)(0.0154)0.2078 0.2660 2.0820 0.1771 0.1789 AUG 2.938 (0.2065)(0.1666)(0.0787)(0.0589)(0.0248)0.0108 0.5102 0.2113 0.2450 0.5669 SEP 4.085 (0.9916)(0.0003)(0.0893)(0.0190)(0.0001)-0.6985 0.3960 0.2287 0.4667 0.7749 OCT 3.073 (0.0866)(0.0020)(0.0433)(0.0001)(0.0001)0.2473 0.3167 0.3009 0.1063 0.4575 1.228 NOV (0.1290)(0.0030)(0.0804)(0.0069)(0.0001)0.3807 0.7723 -0.2623 0.2332 0.4381 DEC 1.585 (0.0001)(0.1987)(0.1065)(0.0084)(0.0001)

Table 5.1. Least Squares Regression Coefficients for Equation (5.1) and Their Significant Levels. The standard deviation, s, for each month is also given.

example, for January the coefficient  $b_1$  is not significant at the 5% significance level ( $\alpha = 0.05$ ) since 0.279 is greater than 0.05, but the R<sup>2</sup> coefficient is significant even at 0.01% significance level ( $\alpha = 0.0001$ ). The significance levels correspond to the "t-test" for the regression coefficients and to the "F-test" for the R<sup>2</sup> coefficients. The standard deviation, s, of the {y} series is also listed since the random component is given by

$$\varepsilon = \sqrt{1 - R^2} s \tag{5.4}$$

as has already been discussed in Chapter 2.

It is interesting to note, that although the multiple regression coefficient R<sup>2</sup> varies for each month from as low as 0.18 to as high as 0.91 it is always significant at the 5% significance level. The months of July and August exhibit the lowest (although significant) correlation coefficients as is expected for Florida. The physical reason for these low correlations is that in the summer most rainfall is convective, whereas in other months there is more cyclonic activity. Rainfall from scattered thunderstorms is simply not as correlated with that of nearby areas as is rainfall from broad cyclonic activity. Thus, on the basis of the regressions shown in Table 5.1, the least squares method would be expected to perform least well in the summer in Florida, but this point is not validated in this thesis.

#### Modified Weighted Average Method (MWA)

For the modified weighted average method the twelve (3x3) covariance matrices of the three index stations have been calculated for each month using equation (2.9) and (2.10), and are shown in Table C.11 (appendix C). Also the monthly standard deviations,  $s_v$ , have been estimated from the known  $\{y\}$  series, and the monthly standard deviations,  $s_y'$  have been calculated by equation (2.11) using the calculated covariance matrices. Notice that although the twelve s values (as calculated from the actual data and which we want to preserve) are different at different percentages of missing values, the twelve  $s'_v$  values (that depend only on the weights a; and the covariance matrix of the index stations) are calculated only once. The correction coefficients f (f =  $s_y/s_y'$ ) for each month and for each different percentage of missing values which must be applied on matrix A (equation 2.21) are shown in Table 5.2.

From this table it can be seen that if the simple weighted average scheme of equation (2.3) were used for the generation, the standard deviation of November would be overestimated (by a factor of approximately 2) and the standard deviation of all other months would be underestimated (e.g., by a factor of approximately 0.5 for the month of January). We also observe that due to small changes of  $s_y$  for different percentages of missing values, the correction factor f does not vary much either, but tends

	(1 - y'y).				
	2%	5%	10%	15%	20%
JAN	1.777	1.777	1.795	1.897	1.872
FEB	1.129	1.142	1.136	1.199	1.188
MAR	1.178	1.207	1.177	1.003	1.009
APR	1.089	0.980	1.061	1.051	1.054
МАҮ	1.269	1.197	1.212	1.222	1.360
JUN	1.214	1.173	1.192	1.228	1.242
JUL	1.338	1.345	1.386	1.390	1.491
AUG	1.424	1.414	1.425	1.432	1.369
SEP	1.313	1.328	1.325	1.210	1.331
OCT	1.258	1.273	1.218	1.229	1.314
NOV	0.533	0.537	0.509	0.583	0.572
DEC	1.161	1.140	1.169	1.172	1.248

Table 5.2. Correction Coefficient, f, for Each Month and for Each Different Percent of Missing Values  $(f = s_y/s'_y)$ .
to be slightly greater the greater the percent of missing values.

The modified weighted average scheme theoretically preserves the mean and variance of the series as has been shown in Chapter 2. But this is true for a series that has been generated by the model and not for a series that is a mix of existing values and values generated (estimated) by This illustrates the difference between the two the model. "generation of data by a model" and "estimation concepts: of missing values by a model." A method for generation of data which is considered "good" in the sense that it preserves first and second order statistics is not necessarily "good" for the estimation of missing values. In fact, it may give statistics comparable to the ones given from a simpler estimation technique which does not preserve the statistics, even as a generation scheme. Theoretically, for a "large" number of missing values, the estimation model operates as a generation model and thus preserves the "desired" statistics, but practically, for this large amount of missing values the "desired" statistics (calculated from the few existing values) are of questionable reliability. Only for augmentation of the time series (extension of the series before the first or after the last point) will the modified weighted average scheme or other schemes that preserve the "desired" statistics be expected to work better than the simple weighted average schemes.

One other disadvantage of the modified weighted average scheme as well as of the least squares scheme is that negative values may be generated by the model. Since all hydrological variables are positive, the negative generated values are set equal to zero, thus altering the statistics of the series. This is also true for all methods that involve a random component and is mainly due to "big" negative values taken on by the random deviate.

The number of negative values, estimated by the MWA method, which have been set equal to zero in the example that follows were 1, 1, 6, 4, and 9 values for the 2%, 5%, 10%, 15% and 20% levels of missing values, respectively.

The effect of the values arbitrarily set to zero cannot be evaluated exactly, but what can be intuitively understood is that a distortion in the distribution is introduced. A transformation that prevents the generation of negative values could be performed on the data before the application of the generation scheme. Such a transformation is, for example, the logarithmic transformation since its inverse applied on a negative value exists, and the mapping of the transformed to the original data and vice versa is one to one (this is not true for the square root transformation).

# Comparison of the MV, RD, NR and MWA Methods

The performance of each method applied for the estimation of the missing values will be evaluated by comparing the estimated series (existing plus estimated

values) to the incomplete series (really available in practice) and to the actual series (unknown in practice, but known in this artificial case). The criteria that will be used for the comparison of the method will be the following:

- (1) the bias in the mean as measured (a) by the difference between the mean of the estimated series,  $\bar{y}_{e}$ , and the mean of the incomplete series,  $\bar{y}_{i}$  (i = 1, 2, 3, 4, 5 for five different percentages of missing values), and (b) by the difference between the mean of the estimated series,  $\bar{y}_{e}$  and the mean of the actual series,  $\bar{y}_{a}$ ;
- (2) the bias in the standard deviation as measured (a) by the ratio of the standard deviation of the estimated series, s<sub>e</sub>, to the standard deviation of the incomplete series, s<sub>i</sub> and (b) by the ratio of the standard deviation of the estimated series, s<sub>e</sub>, to the standard deviation of the actual series, s<sub>a</sub>;
- (3) the bias in the lag-one and lag-two correlation coefficients as measured by the difference of the correlation coefficient of the estimated series,  $r_e$ , to the correlation coefficient of the actual series,  $r_a$ ;
- (4) the bias of the estimation model as given by the mean of the residuals,  $\bar{y}_r$ , i.e., the mean of the differences between the in-filled (estimated) and hidden (actual) values (this is also a check to

detect a consistent over- or under-estimation of the method);

- (5) the accuracy as determined by the variance of the residuals (differences between estimated and actual values) of the whole series,  $s_r^2$ ;
- (6) the accuracy as determined by the variance of the residuals of only the estimated values,  $s_{r.e}^2$ ; and
- (7) the significance of the biases in the mean, standard deviation and correlation coefficients as determined by the appropriate test statistic for each (see appendix A).

Table 5.3 presents the statistics of the actual series (ACT), of the incomplete series (INC) and of the estimated series by the mean value method (MV), by the reciprocal distances method (RD), by the normal ratio method (NR) and by the modified weighted average method (MWA). The mean  $(\bar{y})$ , standard deviation (s), coefficient of variation ( $c_v$ ) coefficient of skewness ( $c_s$ ), lag-one and lag-two correlation coefficients ( $r_1$ ,  $r_2$ ) of the above series considered as a whole have then been calculated.

Regarding comparison of the means, the following can be concluded from Table 5.4:

 the bias in the mean in all cases is not significant at the 5% significance level as shown by the appropriate t-test;

	ÿ	S	cv	cs	r <sub>1</sub>	r <sub>2</sub>
ACT	4.126	3.673	89.040	1.332	0.366	0.134
		2	2% missing	values		
INC	4.116	3.680	89.397	1.346		
MV	4.125	3.663	88.808	1.335	0.371	0.130
RD	4.124	3.674	89.092	1.336	0.367	0.133
NR	4.114	3.666	89.104	1.339	0.368	0.131
MWA	4.113	3.674	89.331	1.342	0.363	0.131
		Ę	5% missing	values		
INC	4.113	3.671	89.249	1.341		
MV	4.101	3.610	88.040	1.352	0.372	0.139
RD	4.127	3.696	89.550	1.359	0.369	0.133
NR	4.105	3.674	89.501	1.349	0.367	0.131
NWA	4.116	3.720	90.386	1.388	0.364	0.126
<del> </del>		10	)% missing •	values		
INC	4.144	3.705	89.405	1.350		
MV	4.134	3.603	87.152	1.346	0.379	0.159

Table 5.3. Statistics of the Actual (ACT), Incomplete (INC) and Estimated Series (MV, RD, NR, MWA).

continued

Table 5.3. Continued.

1 <sup>r</sup> 2
66 0.134
80 0.166
77 0.155
76 0.162
99 0.133
72 0.132
79 0.133
57 0.123
08 0.160
70 0.156
77 0.152
54 0.153

	INC	MV	RD	NR	MWA	
		(	$\bar{y}_{e} - \bar{y}_{i}$ )	g - e ig =, e , e , e		Ī,
28	0.	0.009	0.008	0.002	0.003	4.116
58	0.	-0.012	0.014	-0.008	0.003	4.113
10%	0.	-0.010	0.006	-0.024	-0.017	4.144
15%	0.	-0.089	0.042	0.000	-0.001	4.135
208	0.	0.042	0.149	0.043	0.086	4.082
		(	$\bar{y}_{e} - \bar{y}_{a}$ )			, Ÿa
28	-0.010	-0.001	-0.002	-0.012	-0.013	4.126
5%	-0.013	-0.025	0.001	-0.021	-0.010	
10%	0.018	0.008	0.024	-0.006	0.001	
15%	0.009	-0.020	0.051	0.009	0.008	
20%	-0.044	-0.002	0.105	-0.001	0.042	

- (2) the bias in the mean of the incomplete series is relatively small but becomes larger the higher the percent of missing values;
- (3) at high percents of missing values the NR method gives the less biased mean;
- (4) except for the RD method which consistently overestimates the mean (the bias being larger the higher the percent of missing values), the other methods do not show a consistent over or underestimation.

Regarding comparison of the variances the following can be concluded from Table 5.5:

- (1) Although slight, the bias in the standard deviation is always significant, but this is so because the ratio of variances would have to equal 1.0 exactly to satisfy the F-test (i.e., be unbiased) with as large a number of degrees of freedom as in this study;
- (2) the MV method always gives a reduced variance as compared to the variance of the incomplete series and of the actual series, the bias being larger the higher the percent of missing values;
- (3) the bias in the standard deviation of the incomplete series is small;
- (4) there is no consistent over or under-estimation of the variance by any of the methods (except the MV method);

-	INC	MV	RD	NR	MWA	
			s <sub>e</sub> /s <sub>i</sub>			s <sub>i</sub>
28	1.	0.995	0.998	0.996	0.998	3.680
5%	1.	0.983	1.007	1.001	1.013	3.671
10%	1.	0.972	0.996	0.986	1.005	3.705
15%	1.	0.957	0.988	0.978	0.994	3.671
20%	1.	0.944	1.006	0.973	1.011	3.701
			s <sub>e</sub> /s <sub>a</sub>			sa
28	1.002	0.997	1.000	0.998	1.000	3.673
5%	0.999	0.983	1.006	1.000	1.013	
10%	1.009	0.981	1.004	0.994	1.014	
15%	0.999	0.956	0.988	0.978	0.994	
20%	1.008	0.952	1.014	0.980	1.019	
·						

Table 5.5. Bias in the Standard Deviation

- (5) the MWA method does not give less biased variance even at the higher percent of missing values tested, as compared to the RD and NR methods. Regarding comparison of the correlation coefficients the following can be concluded from Table 5.6:
  - the bias in the correlation coefficients is in all cases not significant at the 5% significance level as shown by the appropriate z-test;
  - (2) the MV method gives the largest bias in the correlation coefficients, the bias increasing the higher the percent of missing values, with a possible effect on the determination of the order of the model;
  - (3) all methods (except the MWA method) consistently overestimate the serial correlation coefficient of the incomplete series but not the serial correlation of the actual series and therefore is not considered a problem;
  - (4) the RD method seems to give a correlogram that closely follows the correlogram of the actual series.

Regarding accuracy of the methods the following can be concluded from Table 5.7:

 no method seems to consistently over or underestimate the missing values at all percent levels, but at high percent levels the missing values are overestimated by all methods;

	INC	MV	RD	NR	MWA	
		(	r <sub>1,e</sub> - r <sub>1,</sub>	a)		r <sub>1,a</sub>
28		0.005	0.001	0.002	-0.003	0.366
5%		0.006	0.003	0.001	-0.002	
10%		0.013	0.014	0.011	0.010	
15%		0.033	0.006	0.013	-0.009	
20%		0.042	0.004	0.011	-0.012	
<u></u>		(	r <sub>2,e</sub> - r <sub>2,</sub>	a)		r <sub>2,a</sub>
28		-0.004	-0.001	-0.003	-0.003	0.134
5%		0.005	-0.001	-0.003	-0.008	
10%		0.025	0.032	0.021	0.028	
15%		-0.001	-0.002	-0.001	-0.011	
20%		0.026	0.022	0.018	0.019	

Table 5.6. Bias in the Lag-One and Lag-Two Correlation Coefficients.

	INC	MV	RD	NR	MWA	
		μ <sub>r</sub> =	Σ(y <sub>e</sub> - y <sub>a</sub> )	/N <sub>O</sub>		No
2%		-0.043	-0.061	-0.570	-0.589	13
5%		-0.440	0.034	-0.380	-0.176	33
10%		0.007	0.156	-0.113	-0.046	62
15%		-0.175	0.338	0.074	0.105	98
20%		0.037	0.502	0.038	0.200	130
		s² r,e	$= \Sigma (y_e -$	y <sub>a</sub> ) <sup>2</sup> /(N <sub>o</sub> -2	)	
2%	<b></b>	5.037	2.874	3.149	4.585	
5%		8.610	3.656	3.411	5.340	
10%		7.892	4.239	3.484	5.187	
15%		7.620	4.630	3.958	5.816	
20%		5.224	4.891	3.681	4.898	

Table 5.7. Accuracy--Mean and Variance of the Residuals N = number of missing values  $N^{\circ} = total number$  of values = 660.

Table 5.7. Continued.

	INC	MV	RD	NR	MWA
		$s_r^2 =$	$\Sigma(y_e - y_a)$	) <sup>2</sup> /(N-2)	
28		0.084	0.048	0.053	0.077
5%		0.406	0.172	0.161	0.252
10%		0.720	0.387	0.318	0.473
15%		1.112	0.675	0.577	0.849
20%		1.016	0.951	0.716	0.953

(2) the NR method is the more accurate method especially at high percents of missing values (i.e., it gives the smaller mean and variance of the residuals).

# Univariate Model

#### Model Fitting

Before considering the problem of missing values the problem of fitting an ARMA(p,q) model to the monthly rainfall series of the south Florida interpolation station will be considered.

The observed rainfall series has been normalized using the square root transformation and the periodicity has been removed by standardization. The reduced series, approximately normal and stationary, is then modeled by an ARMA(p,q) model. The ACF of the reduced series, as shown in Fig. 5.3, implies a white noise process since almost all the autocorrelation coefficients (except at lag-3 and lag-12) lie inside the 95 percent confidence limits.

Of course, it is unsatisfying to accept the white noise process as the "best" model for our series and an attempt is made to fit an ARMA(1,1) model to the series. The selection of an ARMA model and not an AR or MA model is based on the following reasons:

 The observed rainfall series contains important observational errors and so it is assumed to be the sum



Fig. 5.3. Autocorrelation function of the normalized and standardized monthly rainfall series of Station A.

of two series: the "true" series and the observational error series (signal plus noise). Therefore, even if the "true" series obeys an AR process, the addition of the observational error series is likely to produce an ARMA model:

$$AR(p) + white noise = ARMA(p,p)$$

$$AR(p) + AR(q) = ARMA(p+q, max(p,q))$$

$$AR(p) + MA(q) = ARMA(p, p+q)$$

$$(5.5)$$

The same can be said if the "true" series is an MA process and the observational error series an AR process but not if the latter is an MA process or a white noise process:

$$MA(p) + AR(q) = ARMA(q,p+q)$$

$$MA(p) + MA(q) = MA(max(p,q))$$

$$MA(p) + white noise = MA(p)$$
(5.6)

(Granger and Morris, 1976; Box and Jenkins, 1976, Appendix A4.4).

It is understood, that the addition of any observational series to an ARMA process of the "true" series will give again an ARMA process. For example,

$$ARMA(p,q) + white noise = ARMA(p,p) if p > q$$
(5.7)  
= ARMA(p,q) if p  $\leq q$ 

from which it can also be seen that the addition of an observational error may not always change the order of the model of the "true" process.

(2) One other situation that leads exactly, or approximately, to ARMA models is the case of a variable which obeys a simple model such as AR(1) if it were recorded at an interval of K units of time but which is actually observed at an interval of M units (Granger and Morris, 1976, p. 251).

All these results suggest that a number of real data situations are all likely to give rise to ARMA models; therefore, an ARMA(1,1) model will be fitted to the observed monthly rainfall series of the south Florida interpolation station. The preliminary estimate of  $\phi_1$  (equation 3.23) is -0.08163, and the preliminary estimate of  $\theta_1$  (equation 3.21 for k = 0, 1, 2) is the solution of the quadratic equation

$$0.1656 \quad \theta_1^2 + 1.0204 \quad \theta_1 + 0.1656 = 0 \quad . \tag{5.8}$$

Only the one root  $\theta_1 = -0.1667$  is acceptable, the second lying outside the unit circle. These preliminary estimates of  $\phi_1$  and  $\theta_1$  become now the initial values for the determination of the maximum likelihood estimates (MLE). In general, the choice of the starting values of  $\phi$  and  $\theta$  does not significantly affect the parameter estimates (Box and Jenkins, 1976, p. 236), but this was not the case for the



Fig. 5.4. Sum of squares of the residuals,  $\hat{\Sigma}(a_t^2)$ , of an ARMA (1,1) model fitted to the rainfall series of station A.

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<b>.</b>	Initial H	Estimates	Max. Likelihood	Estimates
Model	φ	θ	ф	θ
A	-0.0816	0.0	-0.0088	-0.0989
В	-0.0816	-0.1667	-0.3140	-0.4056
С	0.1	0.0	0.0537	-0.0278
D	-0.4	-0.5	-0.4064	-0.4939

Table 5.8. Initial Estimates and MLE of the Parameters  $\phi$ and  $\theta$  of an ARMA(1,1) model fitted to the rainfall series of station A.

south Florida rainfall series under study. In particular different initial estimates of  $\phi_1$  and  $\theta_1$  have been tested and the MLE of the parameters are compared in Table 5.8. The MLE have been calculated using the IMSL subroutine FTMXL which uses a modified steepest descent algorithm to find the values of  $\phi$  and  $\theta$  that minimize the sum of squares of the residuals (Box and Jenkins, 1976, p. 504).

The drastic changes in parameter values together with the idea that the process may be a white noise process suggest a plot of the sum of squares of the residuals for the visual detection of anomalies. The sum of squares grids and contours are shown in Fig. 5.4. We observe that there is not a well defined point where the sum of squares becomes a minimum but rather a line (contour of the value 641) on which the sum of squares has an almost constant value equal to the minimum. In such case combinations of parameter values give similar sum of squares of residuals and a change in the AR parameter can be nearly compensated by a suitable change in the MA parameter.

From the comparison of the parameters  $\phi$  and  $\theta$ (Table 5.8) of the four ARMA(1,1) models one cannot say that they all correspond to the same process. But this can in fact be illustrated by converting the four models to their "random shock form" (MA( $\infty$ ) processes) or their "invertible form" (AR( $\infty$ ) processes).

An ARMA(1,1) process

$$(1-\phi_1 B) z_t = (1-\theta_1 B) a_t$$
 (5.9)

can be also written as

$$z_t = (1 - \theta_1 B) (1 - \phi_1 B)^{-1} a_t$$
 (5.10)

which can be expanded in the convergent form

$$z_{t} = [1 + (\phi_{1} - \theta_{1})B + \phi_{1}(\phi_{1} - \theta_{1})B^{2} + \phi_{1}^{2}(\phi_{1} - \theta_{1})B^{3} + \dots] a_{t}$$
(5.11)

provided that the stationarity condition ( $|\phi_1| < 1$ ) is satisfied. Then the four models of Table 5.8 become:

(A) : 
$$z_t = a_t + 0.090 a_{t-1} - 0.001 a_{t-2} + \cdots$$
  
(B) :  $z_t = a_t + 0.092 a_{t-1} - 0.029 a_{t-2} + \cdots$  (5.12)  
(C) :  $z_t = a_t + 0.082 a_{t-1} + 0.004 a_{t-2} + \cdots$   
(D) :  $z_t = a_t + 0.088 a_{t-1} - 0.036 a_{t-2} + \cdots$ 

In the same way the ARMA(1,1) model may be written in the "invertible form"

$$(1-\phi_1 B) (1-\theta_1 B)^{-1} z_t = a_t$$
 (5.13)

which can be expanded as

$$[1 - (\phi_1 - \theta_1)B - \theta_1(\phi_1 - \theta_1)B^2 - \theta_1^2(\phi_1 - \theta_1)B^3 - \dots] z_t = a_t$$
(5.14)

given that the invertibility condition ( $|\theta_1| < 1$ ) is satisfied. Then the four models become:

(A) : 
$$z_t = a_t + 0.090 z_{t-1} - 0.009 z_{t-2} + \cdots$$
  
(B) :  $z_t = a_t + 0.092 z_{t-1} - 0.037 z_{t-2} + \cdots$  (5.15)  
(C) :  $z_t = a_t + 0.082 z_{t-1} - 0.002 z_{t-2} + \cdots$   
(D) :  $z_t = a_t + 0.088 z_{t-1} - 0.043 z_{t-2} + \cdots$ 

From the "random shock" form of the four models (equations 5.12) and from their "invertible form" (equations 5.15) the following remarks can be made:

- (1) Although from the comparison of the  $\phi$  and  $\theta$ 
  - coefficients (Table 5.8) of the four ARMA(1,1) models one cannot say that they all correspond to the same process, the comparison of the MA coefficients  $(\theta_1, \theta_2, \theta_3, \ldots)$  of equations (5.12) or the AR coefficients  $(\phi_1, \phi_2, \phi_3, \ldots)$  of equations (5.15) imply that indeed all four models belong to the same process.
- (2) Because the nonzero  $\phi_2$  (and  $\theta_2$ ) coefficients of  $z_{t-2}$ (and  $a_{t-2}$ ) terms while small are of similar magnitude to the coefficients  $\phi_1$  (and  $\theta_1$ ), one cannot say that the "truncated" AR(1) or MA(1) model will fully describe the time series, but instead more terms are needed. On the other hand, we observe that the  $\phi_1$ coefficient so obtained (different for each model) is in the range of 0.082 to 0.090 and is greater than the coefficient  $\phi_1$  that would have been obtained by a direct fitting of an AR (1) model to the series (the latter would be  $\phi_1 = r_1 = 0.0068$ ).
- (3) It should also be noted that all the above models fitted to the series give residuals that pass the portemanteau goodness of fit test. As it can be seen from equation (5.12) the impulse response function (e.g., the weights  $\psi_j$  applied on the  $a_j$ 's when the model is written in the "random shock form") dies off very quickly in all the models, and there is thus no doubt as to the application of the portemanteau test

(see Appendix A). The values of Q for each model (calculated from equation A.1 using K = 60) are:  $Q_A =$ 67.80,  $Q_B = 67.26$ ,  $Q_C = 67.73$  and  $Q_D = 67.39$ , all smaller than the  $\chi^2$  value with 58 degrees of freedom at a 5% significance level,  $\chi^2_{58,5\%} = 79.1$ . It can also be seen that the values of Q for all models are almost equal, suggesting an equally good fit of the series by all the four models.

One other interesting question that could be asked is, given a specific ARMA(p,q) model whether or not this could have arisen from some simpler model. "Simplifications are not always possible as conditions on the coefficients of the ARMA model need to be specified for a simpler model to be realizable" (Granger and Morris, 1976, p. 252). At this stage with coefficients that are so instable it is meaningless to test the four ARMA models for simplification. However, this test will be made after a unique and stable model has been obtained through the following proposed algorithm.

#### Proposed Estimation Algorithm

The problem of estimation of missing values will be combined with the problem of stabilizing the coefficients of the ARMA(1,1) model in a recursive algorithm which will have solved both problems uniquely upon convergence.

The incomplete series  $(\underline{S}_0)$  is filled-in with some initial estimates of the missing values (these initial

estimates can be simply the monthly means or even zeroes as will be shown). Denote by  $\underline{S}_1$  this initial series. An ARMA (1,1) model is fitted to the series  $\underline{S}_1$  and its coefficients  $\phi_1$  and  $\theta_1$  are used to update the first estimates of the missing values. For example, suppose that a gap of size k (k missing values) exists in the series  $\underline{S}_0$ :

Series 
$$\underline{s}_0$$
:  $\ldots$   $z_{t-1}$   $z_t$   $\ldots$   $z_{t+k+1}$   $z_{t+k+2}$   $\ldots$  (5.16)  
Series  $\underline{s}_1$ :  $\ldots$   $z_{t-1}$   $z_t$   $z'_{t+1}$   $\ldots$   $z'_{t+k}$   $z_{t+k+1}$   $z_{t+k+2}$   $\ldots$ 

where  $z'_{t+1}$ , ...,  $z'_{t+k}$  are the initial estimates of the missing values. These values  $z'_{t+1}$ , ...,  $z'_{t+k}$  are then replaced by the <u>forecasted</u> values  $\hat{z}_t(1)$ , ...,  $\hat{z}_t(k)$  by the model, made at origin t and for lead times  $\ell = 1, ..., k$ . These forecasts are the minimum mean square error forward forecasts as developed by Box and Jenkins (1976). For an ARMA(1,1) model with coefficients  $\phi_1$  and  $\theta_1$ , the minimum mean square error forecasts  $\hat{z}_t(\ell)$  of  $z_{t+\ell}$ , where  $\ell$  is the lead time, are:

$$\hat{z}_{t}(\ell) = \phi_{1} z_{t} - \theta_{1} a_{t} , \quad \ell = 1$$

$$\hat{z}_{t}(\ell) = \phi_{1} \hat{z}_{t}(\ell-1) , \quad \ell = 2, \dots, k$$
(5.17)

from which it can be seen that only the one step ahead forecast depends directly on a<sub>t</sub>, and the forecasts at longer lead times are influenced indirectly (Box and Jenkins, 1976, Ch. 5). The forecasting procedure in repeated for the estimation of all the gaps, and the newly estimated values are used in equations (5.17). These forecasts now become the new estimates of the missing values and they replace the old estimates giving the new series  $\underline{S}_2$ . An ARMA(1,1) model is then fitted to the new series  $\underline{S}_2$  and the new coefficients  $\phi_1$  and  $\theta_1$  are found (different from the previous ones). Then the estimated values (forecasts from the previous model) are replaced by the forecasts by the new model, giving the new series  $\underline{S}_3$ , etc. The procedure is repeated until the model and the series stabilize in the sense that the parameters  $\phi_1$  and  $\theta_1$  of the model as well as the estimates of the missing values do not change between successive estimates within a specified tolerance.

Schematically the algorithm is presented in Fig. 5.5 where  $\underline{S}_0$  denotes the incomplete series,  $\underline{M}_0$  the method used for the initial estimation,  $\underline{S}_i$  the estimated series at the ith iteration, and  $\underline{M}_i$  the model (e.g., the set of parameters  $\phi_1$  and  $\theta_1$ ,  $(\phi_1, \theta_1)_i$ ) fitted to the series  $\underline{S}_i$ . The notation  $\underline{M}_i \xrightarrow{\sim} \underline{M}_{i+1}$  and  $\underline{S}_i \xrightarrow{\sim} \underline{S}_{i+1}$  is introduced to denote the stabilization of the model and series respectively after i iterations. The above algorithm will be addressed as RAEMV-U (a recursive algorithm for the estimation of missing values--univariate model).

# Application of the Algorithm on the Monthly Rainfall Series

The proposed recursive algorithm (RAEMV-U) has been applied for the estimation of missing monthly rainfall



Fig. 5.5. Recursive algorithm for the estimation of missing values--univariate model (RAEMV-U). S. denotes the series, and M. the model,  $\overline{(\phi, \theta)}_i$ , at the ith iteration.

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values in the series of the south Florida interpolation station (station 6038). Different levels of percentage of missing values have been tested and the results for the 10% and 20% levels are presented herein. Tables 5.9 and 5.10 show the results for the 10% and 20% levels of missing values respectively. The starting series  $\underline{S}_0$  is the incomplete series (with 10% or 20% the values missing). Four different methods  $\underline{M}_0$  (MV, RD, NR, and zeros) have been applied to the incomplete series,  $\underline{S}_0$ , providing different starting series,  $\underline{S}_1$ , for the algorithm. Thus, its dependence on the initial conditions has also been tested.

# Results of the Method

From Tables 5.9 and 5.10 the following can be concluded:

- (1) The algorithm converges very rapidly and independently of the initial estimates, thus suggesting the convenient replacement of the missing values by zeros to start the algorithm.
- (2) The greater the percent of missing values the slower the algorithm converges (6 iterations were needed for the 10% and 8 for the 20% to obtain accuracy to the third decimal place) as was expected since a larger part of the series is changing its values at each iteration and thus more iterations are needed to achieve equilibrium.

<u>M</u> 0	MV	RD	NR	Zeroes
<u>M</u> 1	0.0255 -0.0208	0.5092 0.5292	0.5018 0.4272	0.5059
<u>M</u> 2	0.4851	0.5010	0.5096	0.4999
	0.4323	0.4166	0.4336	0.4313
<u>M</u> 3	0.5149	0.5110	0.5094	0.5088
	0.4406	0.4355	0.4332	0.4333
$\frac{M}{4}$	0.5087	0.5093	0.5094	0.5094
	0.4322	0.4329	0.4333	0.4333
<u>M</u> 5	0.5096	0.5095	0.5096	0.5095
	0.4335	0.4334	0.4334	0.4334
M <sub>6</sub>	0.5095	0.5095	0.5095	0.5095
	0.4333	0.4333	0.4333	0.4334

Table 5.9. Results of the RAEMV-U Applied at the 10% Level of Missing Values. Upper Value is  $\phi_1$ , Lower Value is  $\theta_1$ .

Table 5.10. Results of the RAEMV-U Applied at the 20% Level of Missing Values. Upper Value is  $\phi_1$ , Lower Value is  $\theta_1$ .

<u>M</u> 0	MV	RD	NR	Zeroes
<u>M</u> 1	0.0954	0.5023	0.5021	0.5756
	-0.0069	0.4173	0.4159	0.2587
<u>M</u> 2	0.0738	0.1167	0.1189	0.2926
	-0.0344	-0.0311	-0.0289	0.1187
<u>M</u> 3	0.0789	0.0369	0.0377	0.0762
	-0.0276	-0.0693	-0.0688	-0.0458
<u>M</u> 4	0.0774	0.0910	0.0908	0.0526
	-0.0296	-0.0125	-0.0128	-0.0503
<u>M</u> 5	0.0778	0.0745	0.0746	0.0863
	-0.0291	-0.0334	0.0333	-0.0184
<u>M</u> 6	0.0777	0.0786	0.0786	0.0756
	-0.0292	-0.0281	-0.0281	-0.0319
<u>M</u> 7	0.0777	0.0775	0.0775	0.0783
	-0.0292	-0.0295	-0.0295	-0.0285
<u>M</u> 8	0.0777	0.0778	0.0778	0.0776
	-0.0292	-0.0291	-0.0291	-0.0293

(3) For a specific percent of missing values the algorithm converges to the same point (e.g., same model and same series) independently of the initial estimates of the missing values.

Diagnostic checking on the residuals from the two final models is performed using the portemanteau goodness of fit test. Denote the two models (at 10% and 20% levels) by  $\underline{M}-\underline{U}^{10}$  and  $\underline{M}-\underline{U}^{20}$  respectively, the U denoting that a univariate model has been fitted to the series. Then

$$\underline{M} - U^{10} : \qquad \phi = 0.5095 \qquad \theta = 0.4333 \qquad (5.18)$$
$$M - U^{20} : \qquad \phi = 0.0777 \qquad \theta = 0.0292.$$

The values of Q for each model are  $Q(M-U^{10}) = 26.54$  and  $Q(M-U^{20}) = 30.22$  (calculated by equation A.1 using K=30) which are both smaller than the  $\chi^2$  value with 28 degrees of freedom at a 5% significance level:  $\chi^2_{28,5\%} = 41.3$ . Notice also that  $Q(M-U^{10}) < Q(M-U^{20})$ , indicating that the final model fitted to the series when 10% of the values were missing has a better fit than the model fitted to the series when 20% of the values were missing as expected.

Also, now that the final ARMA(1,1) model is stable we can ask the question "can it be simplified to an AR(1) plus white noise?". For an ARMA(1,1) process the simplification condition is

$$\frac{1}{1+\phi_{1}^{2}} \ge -\frac{\phi_{1}}{\phi_{1}} \ge 0$$
(5.19)

where

$$\rho_{1} = \frac{\theta_{1}}{\frac{2}{1+\theta_{1}}}$$
(5.20)

(Granger and Morris, 1976, p. 252). For the two models  $\underline{M}-\underline{U}^{10}$  and  $\underline{M}-\underline{U}^{20}$  of equations (5.18) the condition (5.19) gives

$$\underline{M}$$
-U<sup>10</sup> : 0.794 > 0.716 > 0  
M-U<sup>20</sup> : 0.994 > -0.375 ≯ 0 (5.21)

Although the first model barely satisfied the condition for simplification the second model does not, implying that an AR(1) process cannot describe the series as well as an ARMA (1,1) process. This result justified the selection of the ARMA(1,1) model for this rainfall series.

The statistical properties of the two final series (from the 10% and 20% missing values) have also been computed and are shown in Table 5.11 together with the ones of the actual series. The monthly statistics are also shown in Table C.13 (appendix C).

Table 5.11. Statistics of the Actual Series (ACT) and the Two Estimated Series (UN10, UN20).

	ÿ	S	c <sub>v</sub>	cs	r <sub>1</sub>	r <sub>2</sub>
ACT	4.126	3.673	89.04	1.332	0.366	0.134
UN10	4.105	3.609	87.920	1.354	0.384	0.157
UN20	4.043	3.492	86.381	1.373	0.410	0.160

Table 5.12 shows the bias in the mean, standard deviation and lag-one correlation coefficient so that the statistical closeness of the estimated series to the actual one can be evaluated. The bias in the mean and correlation coefficient is not significant at 5% significance level; however, the bias in the standard deviation does not pass the stringent F-test (requiring exact equality of standard deviations) and thus is significant.

	ye-ya	s <sub>e</sub> /s <sub>a</sub>	<sup>r</sup> 1,e <sup>-r</sup> 1,a
UN10	-0.021	0.983	0.018
UN 2 0	-0.083	0.951	0.044

Table 5.12. Bias in the Mean, Standard Deviation and Serial Correlation Coefficient-Univariate Model.

#### Remarks

The forecasting procedure utilized for the estimation 1. is the minimum mean square forward forecasting procedure of Box and Jenkins (1976). Damsleth (1980) introduced the method of optimal between-forecasts, combining the forward forecasts and backforecasts into between-forecasts with a minimum mean square error. He showed that the gain in forecast error by betweenforecasting as compared to forward forecasting (or back-forecasting) an ARMA(1,1) model is proportional to  $|\phi|^{k+1}$  where k is the size of the gap. Thus the gain rapidly becomes small, unless  $|\phi|$  is very close to one and the size of the gap is very small. He also showed that the gain from between-forecasting can be substantial when  $\theta$  is negative. Finally he concluded that "the reduction in forecast error variance by using this between-forecasting method is not very great for stationary series, but may be substantial when the

series is non-stationary" (Damsleth, 1980, p. 39). In our case, the use of the more complicated betweenforecasting procedure does not seem to be justified. It has been shown that the simple Box-Jenkins forecasts work satisfactorily in the sense that rapid convergence to a "statistically acceptable" series occurs.

- 2. It is interesting to note that when the final estimates of the model (parameters of equations 5.18) are provided as initial estimates, the maximum likelihood estimates (calculated by a steepest descent algorithm) are equal to the initial estimates provided. This emphasizes the "uniqueness" of the stable model achieved by the proposed recursive algorithm.
- 3. It will also be interesting to check the threshold level of percent of missing values at which the algorithm starts to diverge. This is expected to happen at some level of percent of missing values (probably greater than 50%) when too much information in the system is changing at each iteration. At such high percents of missing values a more elaborate testing of the final model may also be needed.

## Bivariate Model

# Model Fitting

The lag-one multivariate autoregressive model of equation (4.3), suggested by Matalas (1967), preserves the

lag-zero and lag-one auto- and cross-correlations. When applied to two stations the model is reduced to the bivariate Markov model:

$$\begin{bmatrix} z_{1,t} \\ z_{2,t} \end{bmatrix} = \begin{bmatrix} a_{11} & a_{12} \\ a_{21} & a_{22} \end{bmatrix} \begin{bmatrix} z_{1,t-1} \\ z_{2,t-1} \end{bmatrix} + \begin{bmatrix} b_{11} & 0 \\ b_{21} & b_{22} \end{bmatrix} \begin{bmatrix} \eta_{1,t} \\ \eta_{2,t} \end{bmatrix}$$
(5.22)

where the matrix B is a lower triangular matrix as suggested by Young (1968). The above model has been extensively used for the simultaneous generation of hydrologic series at two sites. An attempt will be made herein, to show how the above model can be used for the estimation of the missing values in one or both of the time series. A recursive algorithm analogous to the one proposed for the univariate case will be presented.

The special case that will be considered is the estimation of the missing values in the series of station 1, given the complete, concurrent, equal length series of station 2.

As has been extensively discussed in Chapter 4 incomplete data sets may result in inconsistent covariance matrices resulting in generated rainfall values that contain complex numbers. Therefore the incomplete series  $\underline{S}_0$  of station 1 is first completed by the use of a simple estimation method  $\underline{M}_0$  (e.g., MV, RD, NR or even replacement of missing values by zeroes) giving the complete series  $\underline{S}_1$ . Denote by <u>S</u> the complete and known series of station 2. Then a bivariate AR(1) model is fitted to the series  $\underline{S}_1$ and  $\underline{S}$ . Actually the model, as in the univariate case, is fitted to the residual series e.g., the normalized and standardized series. The following procedure is followed for the estimation of the parameters (matrices A and B) of the model: The lag-zero and lag-one correlation matrices,  $M_0$  and  $M_1$ , of the residual series are computed

$$M_{0} = \begin{bmatrix} 1 & r_{12}(0) \\ r_{12}(0) & 1 \end{bmatrix}, M_{1} = \begin{bmatrix} r_{11}(1) & r_{12}(1) \\ r_{21}(1) & r_{22}(1) \end{bmatrix} (5.23)$$

Then matrix A is given directly by the multiplication of the matrices  $M_1$  and  $M_0^{-1}$  (equation B.8 of appendix B) and matrix C is computed from equation (B.13). Matrix B is given from the solution of equation  $BB^T = C$ , which in the case of B being a lower triangular matrix reduces to the direct calculation of the elements of B from equations (B.19).

# Proposed Estimation Algorithm

An algorithm analogous to the one for the univariate case is also proposed for the bivariate case. After the incomplete series,  $\underline{S}_0$  has been completed with a simple method  $\underline{M}_0$ , a bivariate AR(1) model is fitted to the complete series  $\underline{S}_1$  and  $\underline{S}$  as described earlier. The parameter matrices A and B of the fitted model  $\underline{M}_1 = (A,B)_1$ , are then used to construct new estimates for the "missing" values in the series  $\underline{S}_1$ . From equation (5.22) we can write that:
$$z_{1,t} = a_{11} z_{1,t-1} + a_{12} z_{2,t-1} + b_{11} \eta_{1,t}$$
 (5.24)

$$z_{2,t} = a_{21} z_{1,t-1} + a_{22} z_{2,t-1} + b_{21} n_{1,t} + b_{22} n_{2,t}$$
 (5.25)

Since the second series is complete and known, equation (5.25) is ignored and only equation (5.24) is considered. Following the Box-Jenkins forecasting procedure, the mean square error forecasts  $\hat{z}_{1,t}(l)$  of  $z_{1,t+l}$ , where l is the lead time, are

$$\hat{z}_{1,t}^{(l)} = a_{11} z_{1,t}^{(l+a_{12} z_{2,t})}, \qquad l = 1$$

$$\hat{z}_{1,t}^{(l)} = a_{11} \hat{z}_{1,t}^{(l-1)} + a_{12} \hat{z}_{2,t}^{(l-1)}, \qquad l = 2, 3, ..., k$$
(5.26)

where k is the number of values missing in each gap. The forecasting procedure is repeated for the estimation of all the gaps always using the newly estimated values in equations (5.26). These estimates then become the new estimates of the missing values, and they replace the old estimates in the series  $\underline{S}_1$  giving the new series  $\underline{S}_2$  and  $\underline{S}$ . Denote this new model by  $\underline{M}_2 = (A,B)_2$ , which is used in the same way as before to update the estimates. The procedure is repeated until convergence occurs in the sense that neither the model  $\underline{M}_i$  nor the series  $\underline{S}_i$  after the ith iteration change between iterations within a specified tolerance ( $\underline{M}_i \xrightarrow{\sim} \underline{M}_{i+1}$  and  $\underline{S}_i \xrightarrow{\sim} \underline{S}_{i+1}$ ).

Schematically the recursive algorithm for the estimation of missing values--bivariate model--1 station to be estimated (RAEMV-B1) is shown in Fig. 5.6.

The algorithm can be generalized to the case where a multivariate model of, say, K stations is used to estimate the missing values of L incomplete stations where  $L \leq K$ . Such a generalized algorithm can be economically written as RAEMV-MK.L. The algorithm for the case of a bivariate model with both records incomplete e.g., two series to be estimated (RAEMV-B2 or in the general form RAEMV-M2.2) is illustrated in Fig. 5.7. The notation is the same as before but two subscripts are used now for the series S, the first denoting the station (1 or 2) and the second denoting the iteration i (i=1, ...). In this case both equations (5.24) and (5.25) would be needed for the estimation of missing values existing in both series.

## Application of the Algorithm on the Monthly Rainfall Series

The case study presented herein involves the estimation of the missing values of the rainfall series of station 6038 using a bivariate AR(1) model with the complete rainfall series of Station 6038. Thus the RAEMV-B1 illustrated in Fig. 5.6 has been used. Again, different levels of percentage of missing values have been tested, and the results for the 10% and 20% missing values are presented in Tables 5.13 and 5.14 respectively. The dependence of the algorithm on the starting values has been tested the same



Fig. 5.6. Recursive algorithm for the estimation of missing values--bivariate model--1 station to be estimated.  $\underline{S}_{i}$  denotes the series, and  $\underline{M}_{i}$  the model,  $(\overline{A}, B)_{i}$ , at the ith iteration.

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Fig. 5.7. Recursive algorithm for the estimation of missing values--bivariate model--2 stations to be estimated (RAEMV-B2).  $S_{1,i}$  ( $S_{2,i}$ ), denotes the series of station 1'(station 2), and  $\underline{M}_{i}$  the model, (A,B)<sub>i</sub>, at the ith iteration.

way as for the univariate case, e.g., by providing different initial series estimated by four different methods  $\underline{M}_0$  (MV, RD, NR and zeroes).

Tables 5.13 and 5.14 show the cross-correlation matrices  $M_0$  and  $M_1$  at each iteration i and the model  $\underline{M_i} = (A,B)_i$ . It is interesting to follow the changes of the cross-correlation coefficients at each time step. Also notice that the autocorrelation coefficient (see equation 5.23) of the first series changes at each iteration (since new estimates of the missing values replace the old ones) but the autocorrelation coefficient of the second series remains unchanged (since the second series is complete and known).

From Tables 5.13 and 5.14 the following similar conclusions to the univariate case can be drawn:

- (1) The algorithm converges rapidly, independently of the starting point (initial series). Thus, initial estimation of the missing values is not needed, and they may as well be replaced by zeroes.
- (2) The convergence seems to be less sensitive to the percent of values missing, since in both the 10% and 20% levels convergence has been achieved in three to four iterations.
- (3) For a specific percent of missing values the algorithm converges to the same point (e.g., same model, same series, and same correlation matrices) independently of the initial estimates of the missing values.

i	M		M <sub>1</sub>		A	<u></u> i	E	3
				M	$I_0 = MV$			
_	1.	0.330	0.004	0.137	-0.046	0.152	0.990	0.
1	0.330	1.	0.042	0.315	-0.070	0.338	0.286	0.902
	1.	-0.005	0.038	0.194	0.039	0.194	0.980	0.
2	-0.005	1.	0.065	0.315	0.067	0.316	-0.071	0.944
_	1.	0.025	0.049	0.202	0.044	0.201	0.978	0.
3	0.025	1.	0.069	0.315	0.061	0.314	-0.043	0.946
	1.	0.025	0.049	0.201	0.044	0.200	0.979	0.
4	0.025	1.	0.068	0.315	0.061	0.314	-0.042	0.946
				<u>I</u>	$l_0 = RD$			
1	1.	0.554	0.124	0.249	-0.021	0.261	0.968	0.
Ŧ	0.554	1.	0.201	0.315	0.038	0.294	0.492	0.811
2	1.	0.026	0.042	0.196	0.037	0.195	0.980	0.
Ζ	0.026	1.	0.070	0.315	0.062	0.314	-0.039	0.946
2	1.	0.025	0.048	0.201	0.043	0.200	0,979	0.
3	0.025	1.	0.069	0.315	0.061	0.314	-0.042	0.946
	1.	0.025	0.049	0.201	0.044	0.200	0.979	0.
4	0.025	1.	0.068	0.315	0.061	0.314	-0.042	0.946
				Conti	nued			

Table 5.13. Results of the RAEMV-B1 Applied at the 10% Level of Missing Values.

Table 5.13. Continued.

				M	$I_0 = NR$			
1	1.	0.543	0.126	0.261	-0.022	0.273	0.965	0.
T	0.543	1.	0.187	0.315	0.022	0.303	0.478	0.819
•	1.	-0.002	0.046	0.199	0.046	0.199	0.979	0.
۷	-0.002	1.	0.069	0.315	0.070	0.316	-0.070	0.944
2	1.	0.026	0.050	0.203	0.045	0.201	0.978	0.
5	0.026	1.	0.069	0.315	0.061	0.314	-0.042	0.946
Λ	1.	0.025	0.049	0.202	0.044	0.200	0.978	0.
4	0.025	1.	0.068	0.315	0.061	0.314	-0.042	0.946
				M	<sub>0</sub> = zero	pes		
1	1.	0.258	0.463	0.172	0.448	0.057	0.885	0.
Т	0.258	1.	0.048	0.315	-0.036	10.385	0.247	0.915
S	1.	0.042	0.061	0.225	0.059	0.222	0.973	0.
۷	0.042	1.	0.081	0.315	0.068	0.313	-0.033	0.946
S	1.	0.029	0.048	0.203	0.042	0.201	0.978	0.
2	0.029	1.	0.070	0.315	0.061	0.314	-0.038	0.946
Λ	1.	0.025	0.049	0.201	0.043	0.200	0.979	0.
4	0.025	1.	0.068	0.315	0.061	0.314	-0.042	0.946

						<u>M</u> ;		
i	<sup>M</sup> 0		<sup>M</sup> 1		A		E	5
				M	0 = MV			
_	1.	0.523	0.342	0.251	0.290	0.100	0.936	0.
T	0.523	1.	0.257	0.315	0.126	0.249	0.446	0.831
-	1.	-0.025	0.369	0.307	0.377	0.316	0.874	0.
2	-0.025	1.	0.256	0.315	0.264	0.322	-0.253	0.876
n	1.	-0.023	0.389	0.333	0.393	0.337	0.857	0.
3	-0.012	1.	0.253	0.315	0.257	0.319	-0.255	0.877
	1.	-0.012	0.389	0.332	0.393	0.337	0.858	0.
4	-0.012	1.	0.253	0.315	0.257	0.319	-0.254	0.877
				<u>M</u>	0 = RD			
1	1.	0.588	0.320	0.290	0.228	0.156	0.939	0.
т	0.588	1.	0.262	0.315	0.117	0.246	0.510	0.795
с С	1.	-0.012	0.368	0.315	0.375	0.383	0.872	0.
2	-0.023	1.	0.257	0.315	0.264	0.321	-0.254	0.875
C	1.	-0.012	0.388	0.334	0.392	0.338	0.857	0.
3	-0.012	1.	0.253	0.315	0.257	0.319	-0.255	0.877
٨	1.	-0.012	0.388	0.333	0.393	0.337	0.858	0.
4	-0.012	1.	0.253	0.315	0.257	0.318	-0.254	0.877
				Conti	nued			

Table 5.14. Results of the RAEMV-B Applied at the 20% Level of Missing Values.

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				M	$_0 = NR$			
-	1.	0.611	0.324	0.273	0.252	0.119	0.941	0.
T	0.611	1.	0.279	0.315	0.137	0.232	0.534	0.777
0	1.	-0.022	0.372	0.311	0.379	0.320	0.872	0.
2	-0.022	1.	0.258	0.315	0.265	0.321	-0.253	0.875
2	1.	-0.012	0.389	0.333	0.393	0.338	0.857	0.
د	-0.012	1.	0.253	0.315	0.257	0.319	-0.255	0.877
Λ	1.	-0.012	0.389	0.332	0.393	0.337	0.857	0.
4	-0.012	1.	0.253	0.315	0.257	0.318	-0.254	0.877
				M	n_= zero	es		
-	1.	0.321	0.601	0.201	0.599	0.009	0.799	0.
T	0.321	1.	0.195	0.315	0.104	0.282	0.253	0.909
n	1.	0.006	0.423	0.340	0.421	0.337	0.841	0.
Ζ	0.006	1.	0.228	0.315	0.226	0.314	-0.233	0.892
n	1.	-0.012	0.392	0.332	0.397	0.337	0.856	0.
3	-0.013	1.	0.249	0.315	0.253	0.319	-0.255	0.878
л	1.	-0.013	0.390	0.333	0.394	0.338	0.857	0.
4		-	0 050	0 215	0 0 5 7	0 210		0 077
	-0.013	⊥.	0.253	0.315	0.25/	0.319	-0.255	0.8//

(4) For a different percent of missing values the same series converges to a "different" point, but this is reasonable and expected since the constant information (existing values in the series) is different in each case, and a different model thus describes it better.

The statistical properties of the two final series (from the 10% and 20% missing values) are shown in Table 5.15 together with the ones of the actual series. The monthly statistics are also shown in Table C.14 (appendix C). Table 5.16 shows the statistical closeness of the two estimated series to the actual one. Again, the bias in the mean and correlation coefficient is not significant at the 5% significance level, but the bias in the standard deviation is.

·····						
	Ŷ	s	°r	cs	r <sub>l</sub>	r <sub>2</sub>
ACT	4.126	3.673	89.04	1.332	0.366	0.134
B10	4.096	3.610	88.132	1.358	0.382	0.162
B20	4.077	3.523	86.421	1.341	0.416	0.165

Table 5.15. Statistics of the Actual Series (ACT) and the Two Estimated Series (B10 and B20).

	y <sub>e</sub> -y <sub>a</sub>	s <sub>e</sub> /s <sub>a</sub>	r <sub>1,e</sub> -r <sub>1,a</sub>
B10	-0.030	0.983	0.016
B20	-0.049	0.959	0.050

Table 5.16. Bias in the Mean, Standard Deviation and Serial Correlation Coefficient-Bivariate Model.

## CHAPTER 6

## CONCLUSIONS AND RECOMMENDATIONS

#### Summary and Conclusions

The objective of this study was to compare and evaluate different methods for the estimation of missing observations in monthly rainfall series. The estimation methods studied reflect three basic ideas:

- (1) the use of regional information in four simple techniques:
  - mean value method (MV),
  - reciprocal distance method (RD),
  - normal ratio method (NR),
  - modified weighted average method (MWA);
- (2) the use of a univariate stochastic (ARMA) model that describes the time correlation of the series;
- (3) the use of a multivariate stochastic (ARMA) model that describes the time and space correlation of the series.

An algorithm for the recursive estimation of the missing values in a time series using the fitted univariate or multivariate ARMA model has been proposed and demonstrated. Apparently, the idea of the recursive estimation of missing values is known (Orchard and Woodbury, 1972; Beale and

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Little, 1974), as well as the idea of using the fitted model to directly derive the estimates (Brubacher and Wilson, 1976; Damsleth, 1979). However it appears that a method which combines the above two ideas simultaneously in a recursive estimation of the missing values with parallel updating of the model has not been used before.

The proposed algorithm is general and can be used for the estimation of the missing values in any series that can be described by an ARMA model.

On the basis of the data from the four south Florida rainfall stations used in the analysis, the following conclusions can be drawn:

- (1) All the simplified estimation techniques give unbiased (overall and monthly) means and correlation coefficients at the 5% significance level even for as high as 20% missing values.
- (2) At high percentages of missing values (greater than 10%) the MV method gives the more biased (although not significantly so) correlation coefficients.
- (3) All methods give a slightly biased overall variance but unbiased monthly variance at the 5% significance level, and the MV method gives the most biased variances for all percentages of missing values.
- (4) The NR method gives the most and the MV the least accurate estimates, at almost all levels of percent missing values.

- (5) The proposed recursive algorithm works satisfactorily in both the univariate and bivariate case. It converges rapidly and independently of the initial estimates and gives unbiased means and correlation coefficients at the 5% significance level.
- (6) The use of a bivariate model as compared to a univariate one did not improve the estimates except for a slight improvement at 20% missing values. However, the use of a multivariate model based on three or four nearby stations is expected to give much better estimates. The use of three adjacent stations is the main reason for the better performance of the NR method over the more sophisticated univariate and bivariate ARMA models which use only zero and one additional stations.

If the purpose of estimation is to calculate the historical statistics of the series (e.g., mean, standard deviation, and autocorrelations) the selection of the method matters little, and the simplest one may be chosen. However, if it is desired to fit an ARMA model to the incomplete series, to be used, say, to construct forecasts, the estimation of the missing values and the parameters of the model by the proposed recursive algorithm is recommended. In this case the equilibrium state (i.e., final series and parameters of the model) achieved upon convergence is unique, depending only on the existing information in the

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system (available data) and not on any external information added to the system (by the replacement of the missing values with some derived estimates). The only assumption made is that the <u>order</u> of the ARMA model to be fitted to the series is known. In practical situations this is seldom a problem since the latter can be determined from the complete part of the series or from a series with similar characteristics. For example, if an ARMA(1,1) model is known to fit the monthly rainfall series well at a couple of nearby stations, there is little doubt that it will fit the incomplete monthly rainfall series equally well at the station of interest. Upon convergence, the recursive algorithm then gives the "best" estimates of the parameters of the model.

#### Further Research

Further research should include:

- application of the simple estimation techniques in short records where the biases may be significant for the methods with the poorer performance;
- (2) test of the sensitivity of the recursive algorithm to the selection of the model (order of the model) when more than one model fits the data equally well;
- (3) derivation of the threshold percent of missing values after which the algorithm diverges;
- (4) application to the estimation of missing values in other hydrological series, e.g., runoff;

- (5) trials of different forecasting procedures and determination of improvements obtained by the "betweenforecasting procedure" in cases of a large number of single-value gaps, e.g., use of the average of a backwards and forwards ARMA model forecast;
- (6) application of the concept of "missing values" for the estimation of erroneous values or outliers in a series to avoid errors when using the data, say, to construct forecasts; and
- (7) estimation of values in a series that are affected by unusual circumstances, thereby permitting a measure of the magnitude of the unusual circumstance and the estimation of the effect of similar circumstances in the future (e.g., effect of a drought on water supply).

#### APPENDIX A

#### DEFINITIONS

### 1. Strict stationarity

A stochastic process is said to be strictly stationary if its statistics (e.g., mean, variance, serial correlation) are not affected by a shift in the time origin, that is, if the joint probability distribution associated with n observations  $(z_1, z_2, \ldots, z_n)_t$  made at time origin t, is the same as that associated with n observations  $(z_1, z_2, \ldots, z_n)_{t+k}$  made at time origin t+k. In other words, z(t) is a strictly stationary process when the two processes z(t) and z(t+k) have the same statistics for any k.

#### 2. Weak stationarity

Weak stationarity of order f is when the moments of the process up to an order f depend only on time differences. Usually by weak stationarity we refer to second order stationarity, e.g., fixed mean and an autocovariance matrix that depends only on time differences (i.e., lags).

#### 3. Gaussian process

If the probability distribution associated with any set of times is a multivariate normal distribution, the process

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is called a normal or Gaussian process. Since the multivariate normal distribution is fully described by its first and second order moments it follows that weak stationarity and an assumption of normality imply strict stationarity.

#### 4. Non-stationarity

A stochastic process is said to be nonstationary if its statistical characteristics change with time. A homogeneous nonstationary process of order d is a process, for which the dth difference  $\nabla^d z_t$  is a stationary process. For example a first order homogeneous nonstationary process is one that exhibits homogeneity apart from constant (e.g., a linear trend), and a second order nonstationary is the one that exhibits homogeneity apart from constant and slope (e.g., a parabolic trend).

### 5. Circular stationarity

A stochastic process is said to be circularly stationary with period  $\tau$ , if the multivariate probability distribution of  $\tau$  observations  $(z_1, z_2, \ldots, z_{\tau})_t$  made at time origin t, is the same as that associated with  $\tau$ observations  $(z_1, z_2, \ldots, z_{\tau})_{t+\tau k}$  made at time origin  $t + \tau k$ , for  $k = 1, 2, \ldots$ . For example, a monthly hydrologic series has a period of 12 months, i.e.,  $\tau = 12$  and circular stationarity suggests that the probability distribution of a value of a particular month is the same for all the years.

# 6. Stationarity condition

A linear process can be always written in the random shock form:

$$\tilde{z}_{t}^{\circ} = \psi(B) a_{t}^{\circ}$$
 (A.1)

where B is the backward shift operator defined by  $Bz_t = z_{t-1}$ ; hence  $B^m z_t = z_{t-m}$  and

$$\psi(B) = 1 + \psi_1 B + \psi_2 B^2 + \dots$$
 (A.2)

is the so called transfer function of the linear system and is the generating function of the  $\psi$  weights. For the process to be stationary the  $\psi$  weights must satisfy the condition that  $\psi(B)$  converges on or within the unit circle, e.g., for all  $|B| \leq 1$ .

#### 7. Invertibility condition

The above model may also be written in the inverted form

$$\psi^{-1}(B) \hat{z}_{t} = a_{t}$$
 (A.3)

or

$$\pi(B) \quad \tilde{z}_{t} = a_{t} \tag{A.4}$$

where

$$\pi$$
 (B) = 1 -  $\pi_1 B$  -  $\pi_2 B^2$  - . . (A.5)

is the generating function of the  $\pi$  weights. For the process to be invertible the  $\pi$  weights must satisfy the condition that  $\pi(B)$  converges for all  $|B| \leq 1$ , that is on or within the unit circle. The invertibility condition is independent of the stationarity condition and is applicable also to the nonstationary linear models. The requirement of invertibility is needed in order to associate the present values of the process to the past values in a reasonable manner, as will be shown below.

## 8. Duality between AR and MA processes

In a stationary AR(p) process,  $a_{t}$  can be represented as a finite weighted sum of previous  $\overset{\sim}{z}$  's,

$$\phi(B) \quad z_t = a_t \tag{A.6}$$

or  $z_t^{\sim}$  as an infinite weighted sum of previous a's

$$\hat{z}_{t}^{\circ} = \phi^{-1}(B) a_{t}$$
 (A.7)

Also, in an invertible MA(q) process,  $z_t^{\circ}$  can be represented as a finite weighted sum of previous a's,

$$\hat{z}_{t}^{\circ} = \theta(B) a_{t}^{\circ}$$
 (A.8)

or  $\textbf{a}_{+}$  as an infinite weighted sum of previous  $\overset{\sim}{z}\textbf{'s}$ 

$$\theta^{-1}(B) \quad \tilde{z}_{t} = a_{t} \tag{A.9}$$

In other words, a finite AR process is equivalent to an infinite MA process, and a finite MA process to an infinite AR process. This principle of duality has further aspects, e.g., there is an inverse relationship between the autocorrelation and partial autocorrelation functions of AR and MA processes.

# 9. <u>Physical interpretation of stationarity and</u> <u>invertibility</u>

Consider an AR(1) process  $(1 - \phi_1 B) z_t = a_t$ . For this process to be stationary, the root of the polynomial  $1 - \phi_1 B = 0$  must lie outside the unit circle, which implies that  $B = \phi_1^{-1}$  must be greater than one, or  $|\phi_1| < 1$ . The process can be also written

$$z_{t} = \phi_{1}z_{t-1} + a_{t}$$

$$z_{t+1} = \phi_{1}^{2}z_{t-1} + \phi_{1}a_{t} + a_{t+1}$$

$$z_{t+2} = \phi_{1}^{3}z_{t-1} + \phi_{1}a_{t} + \phi_{1}a_{t+1} + a_{t+2} \text{ etc.}$$
(A.10)

When  $|\phi_1| > 1$  (or  $|\phi_1| = 1$ ) the effect of the past on the present value of the time series increases (or stays the

same) as the series moves into the future. Only when  $|\phi_1| < 1$  (stationary process) does the effect of the past on the present decrease the further we move into the past, which is a reasonable and acceptable hydrologic fact (Delleur and Kavvas, 1978).

Consider now an MA(1) process  $z_t = (1 - \theta_1 B)a_t$ . The invertibility condition implies that  $|\theta_1| < 1$ . The process can also be written in the form:

$$a_{t} = \frac{1}{1 - \theta_{1}B} z_{t}$$
(A.11)

where the polynomial  $(1-\theta_1 B)^{-1}$  can be expanded in an infinite sum of convergent series only if  $|\theta_1| < 1$ . To illustrate the need for invertibility let us assume that  $|\theta_1| > 1$ . Then (A.11) can be written as

$$a_{t} = -\frac{1}{\theta_{1}B} \frac{1}{(1 - \frac{1}{\theta_{1}B})} z_{t}$$
 (A.12)

and since  $\left|\frac{1}{\theta_1 B}\right| < 1$ , it can be expanded to the form

$$a_{t} = -\left(\frac{1}{\theta_{1}B} + \frac{1}{\theta_{1}^{2}B^{2}} + \frac{1}{\theta_{1}^{3}B^{3}} + \dots\right) z_{t}$$
 (A.13)

or

$$a_{t} = -\frac{1}{\theta_{1}} z_{t+1} - \frac{1}{\theta_{1}^{2}} z_{t+2} - \frac{1}{\theta_{1}^{3}} z_{t+3} - \dots$$
 (A.14)

which implies that future values are used to generate the present values. It becomes clear that the invertibility condition is required in order to assure hydrologic realizability.

## 10. The portemanteau lack of fit test

The portemanteau lack of fit test (Box and Jenkins, 1976, Ch. 8) considers the first K autocorrelations  $r_k(\hat{a})$ ,  $k = 1, 2, \ldots, K$ , of the fitted residual series  $\hat{a}$  of an ARIMA(p,d,q) process, to detect inadequacy of the model. It can be shown (Box and Pierce, 1970) that, if the fitted model is appropriate,

$$Q = (N-d) \sum_{k=1}^{K} r_k^2 (\hat{a})$$
 (A.15)

is approximately distributed as  $\chi^2(K-p-q)$  where K-p-q is the number of degrees of freedom, N is the total length of the series, and (N-d) is the number of observations used to fit the model. The adequacy of the model may be checked by comparing Q with the theoretical chi-square value  $\chi^2(K-p-q)$  of a given significance level. If Q <  $\chi^2(K-p-q)$ ,  $a_t$  is an independent series and so the model is adequate, otherwise the model is inadequate.

For the choice of K, Box and Jenkins suggest it to be "sufficiently large so that the weights  $\psi_j$  in the model, written in the form

$$\widetilde{\mathbf{w}}_{t} = \phi^{-1}(\mathbf{B}) \quad \theta(\mathbf{B}) \quad \mathbf{a}_{t} = \psi(\mathbf{B}) \quad \mathbf{a}_{t}$$
(A.16)

will be negligibly small after j = K" (Box and Jenkins, 1976, p. 221). The IMSL subroutine FTCMP (IMSL - 0007, Ch. F) uses a value of K equal to N/10 + p + q to perform the portemanteau test.

Ozaki (1977) points out that "for the application of the portemanteau test, fast dying off of the impulse response function (weights  $\psi_j$ ) of the model is a necessary condition" (Ozaki, 1977, p. 298). In cases where the impulse response function dies off rather slowly (possibly due to the near-nonstationarity of the model) when compared with the length of the series, the applicability of the portemanteau test is doubtful since the autocorrelations of the residuals may not be reliable at large lags.

#### 11. Cumulative periodogram test

Another method used in the diagnostic checking stage of the Box-Jenkins procedure is the cumulative periodogram checking of the residuals. The normalized (area under the curve equal to one) cumulative periodogram for frequencies, f, between 0 and 0.5, of the fitted residuals  $\hat{a_t}$ , is compared with the theoretical cumulative periodogram of a white noise series which is a straight line joining the points (0, 0) and (0.5, 1). A periodicity in the residuals at frequency  $f_i$  is expected to show up as a deviation from the straight line at this frequency. Kolmogorov-Smirnov probability limits can be drawn on the cumulative periodogram plot to test the significance of such deviations. For a given level of significance  $\alpha$ , the limit lines are drawn at distances  $\pm K_{\alpha}/N'$  above and below the theoretical straight line, where N' = (N-2)/2 for N even and N' = (N-1)/2 for N odd. Approximate values of  $K_{\alpha}$  for different levels of significance  $\alpha$ , are:

 $\alpha$  0.01 0.05 0.10 0.20 0.25  $K_{\alpha}$  1.63 1.36 1.22 1.07 1.02 (Box and Jenkins, 1976, p. 297). So, if more than  $\alpha$ N of the plotted points fall outside the probability lines, the residual series may still have some periodicity; otherwise it may be concluded that the residuals are independent.

In practice, "because the a's are fitted values and not the true a's, we know that even when the model is correct they will not precisely follow a white noise process" and thus the cumulative periodogram test provides only a "rough guide" to the model inadequacy checking (Box and Jenkins, 1976, p. 297). 12. Akaike Information Criterion (AIC)

The AIC for an ARMA(p,q) model is given by

AIC(p,q) = N 
$$\log(\sigma_a^2)$$
 + 2(p+q+2) + Nlog2  $\pi$  + N

where  $\hat{\sigma}_a^2$  is the MLE of the residual variance given by

$$\hat{\sigma}_{a}^{2} = \frac{1}{N-p-q} S(\underline{\phi}, \underline{\theta})$$
(A.17)

and  $\underline{\phi}$ ,  $\underline{\theta}$  are the vectors of the parameters  $\phi$ ,  $\theta$  which minimize the sum of squares of the residuals  $\hat{a}_{+}$ 

$$S(\underline{\phi},\underline{\theta}) = \sum_{t=1}^{N} (\hat{a}_{t})^{2}$$
 (A.18)

For the purpose of comparison of models the definition of AIC can be replaced by

AIC(p,q) = N 
$$\log(\hat{\sigma}_{a}^{2}) + 2(p+q)$$
 . (A.19)

Ozaki (1977) demonstrates that the inherent difficulties associated with the Box-Jenkins procedure (identification, estimation and diagnostic checking) for the selection of the model, when several models fit the data equally well, can be overcome by using the MAICE (minimum AIC estimation) procedure as the only objective criterion for the selection of the "best" approximating model among a set of possible models. He also points out that the AIC "measures both the fit of a model and the unreliability of a model" (Ozaki, 1977, p. 290).

# 13. Positive definite (semidefinite) matrix

A real symmetric matrix A is called positive definite (semidefinite) if and only if

$$X^{T} A X > 0$$
 (>0) (A.20)

for all vectors  $X \neq 0$ . The two following theorems hold:

<u>Theorem 1</u>: A matrix A is positive (semi-) definite if and only if all its characteristic values (i.e., eigenvalues) are (non-negative) positive.

<u>Theorem 2</u>: A matrix A is positive (semi-) definite if and only if all the successive principal minors of A are (non-negative) positive.

An obvious corollary of the above is that a positive semidefinite matrix is positive definite if and only if it is nonsingular i.e., none of its characteristic values are zero (Gantmacher, 1977, p. 305).

# 14. Test for differences in the means of two normal populations

Let  $\mu_1$ ,  $\mu_2$  denote the population means of two normal distributions and  $\bar{x}_1$ ,  $\bar{x}_2$  the sample means respectively. Let also assume that the variance of the two normal distributions are equal but unknown. The hypothesis  $H_0: \mu_1 = \mu_2$  versus  $H_a: \mu_1 \neq \mu_2$  is tested by calculating the statistic

$$t = \frac{\bar{x}_1 - \bar{x}_2}{s\sqrt{1/N_1 + 1/N_2}}$$
(A.21)

where

$$s^{2} = \frac{(N_{1}-1) s_{1}^{2} + (N_{2}-1) s_{2}^{2}}{N_{1} + N_{2} - 2}$$
(A.22)

which has a t distribution with  $N_1 + N_2 - 2$  degrees of freedom. The H<sub>o</sub> is rejected if

$$|t| > t_{1-\alpha/2} ; N_1 + N_2 - 2$$
 (A.23)

Although the test is based on sample normality, for large samples, the Central Limit Theorem enables us to use the test as approximate test for nonnormal samples. If the two populations are of equal length,  $N_1 = N_2 = N$ , then equation (A.21) reduces to

$$t = \frac{\bar{x}_1 - \bar{x}_2}{\sqrt{(s_1^2 + s_2^2)/N}}$$
(A.24)

# 15. <u>Test for equality of variances of two normal</u> <u>distributions</u>

Let  $\sigma_1^2$ ,  $\sigma_2^2$  denote the population variances and  $s_1^2$ ,  $s_2^2$  the sample variances of two normal distributions. The hypothesis  $H_0: \sigma_1^2 = \sigma_2^2$  versus  $H_a: \sigma_1^2 \neq \sigma_2^2$  is tested by calculating the statistic

$$F_{c} = s_{1}^{2} / s_{2}^{2}$$
 (A.25)

where  $s_1^2$  is the larger sample variance.  $F_c$  is distributed as an F distribution with  $N_1 - 1$  and  $N_2 - 1$  degrees of freedom where  $N_1$  is the length of the sample having the larger variance and  $N_2$  is the length the sample with the smaller variance.  $H_o$  is rejected if

$$F_{C} > F_{N_{2}-1}^{N_{1}-1}; 1 - \alpha$$
 (A.25)

## 16. Test for equality of correlation coefficients

Let  $\rho$  denote the population correlation coefficient and r the sample estimate of  $\rho$ . If the sample size is moderately large (N > 25) then the quantity W is approximately normally distributed with mean  $\$  and variance 1/N-3 where

$$W = \frac{1}{2} \ln \left( \frac{1+r}{1-r} \right)$$
 (A.27)

and

$$\omega = \frac{1}{2} \ln \left( \frac{1 + \rho}{1 - \rho} \right) .$$
 (A.28)

To test the hypothesis  $H_0: \rho = r$  against the alternative  $H_a: \rho \neq r$  the quantity

$$z = (W - \omega) \sqrt{N-3}$$
 (A.29)

can be considered to be normally distributed with zero mean and unit variance. If  $|z| > z_{1-\alpha/2}$  (z is the standard normal variable), H<sub>o</sub> is rejected (see Haan, 1977, p. 223).

#### APPENDIX B

### DETERMINATION OF MATRICES A AND B OF THE MULTIVARIATE AR(1) MODEL

Determination of matrix A

The multivariate lag-one autoregressive model is written as

$$Z_{t} = A Z_{t-1} + B N_{t}$$
(B.1)

Post-multiplying both sides of equation (B.1) by  $Z_{t-1}^{T}$  and taking expectations it becomes:

$$E[Z_{t}Z_{t-1}^{T}] = A E[Z_{t-1}Z_{t-1}^{T}] + B E[N_{t}Z_{t-1}^{T}]$$
(B.2)

By definition

$$M_0 = E[Z_t Z_t^T]$$
(B.3)

and

$$M_{1} = E[Z_{t}Z_{t-1}^{T}]$$
(B.4)

and from the assumption of weak stationarity

$$M_0 = E[Z_{t-1}Z_{t-1}^T] . (B.5)$$

Also from the independent uncorrelated process  ${\tt N}_{\tt t}$ 

$$E[N_{t}Z_{t-1}^{T}] = 0$$
 (B.6)

so that equation (B.2) becomes

$$M_1 = A M_0 \tag{B.7}$$

and solving for the parameter matrix A

$$A = M_1 M_0^{-1}$$
 (B.8)

Determination of matrix B

Post-multiplying equation (B.1) by  $Z_t^T$  and taking expectations in both sides it becomes

$$E[Z_{t}Z_{t}^{T}] = A E[Z_{t-1}Z_{t}^{T}] + B E[N_{t}Z_{t}^{T}]$$

$$= A E[Z_{t}Z_{t-1}^{T}]^{T} + B E[N_{t}(AZ_{t-1} + BN_{t})^{T}]$$
(B.9)
$$= A E[Z_{t}Z_{t-1}^{T}]^{T} + B E[N_{t}Z_{t-1}^{T}] A^{T} + B E[N_{t}N_{t}^{T}] B^{T}$$

Because  $E[N_tN_t^T] = I$ , an identity matrix, and  $E[N_tZ_{t-1}^T] = 0$ equation (B.9) can be written

$$M_0 = A M_1^T + B B^T$$
 (B.10)

By substituting A from equation (B.8) and solving for B  $\boldsymbol{B}^{\mathrm{T}}$ 

$$B B^{T} = M_{0} - M_{1} M_{0}^{-1} M_{1}^{T}$$
(B.11)

Solution of equation B  $B^{T} = C$ 

The right hand side of equation (B.11) involves the lag-zero and lag-one correlation matrices which can be estimated from the historical data and thus is a known guantity C. The problem that remains now, is to solve equation

$$B B^{T} = C$$
 (B.12)

for B. A necessary and sufficient condition to have a real solution for B is that C must be a positive semidefinite matrix.

It can be proven (Valencia and Schaake, 1973) that if the correlation matrices  $M_0$  and  $M_1$  have been calculated using equal length records for all m sites, then the matrix

$$C = M_0 - M_1 M_0^{-1} M_1^{T}$$
(B.13)

is always positive semidefinite and so a real solution for the matrix B exists. But this solution for B is not unique. An infinite number of matrices B exist that satisfy (B.12).

<u>Proof</u>: Let B denote a matrix solution of equation (B.12) and K denote an (mxm) matrix such that K  $K^{T} = I$  where I is an (mxm) identity matrix. A matrix  $B_{0}$  defined as

$$B_0 = B K \tag{B.14}$$

may be used in place of B in equation (B.12) since

$$B_0 B_0^T = B K (B K)^T = B K K^T B^T = B B^T$$
 (B.15)

There exists more than one matrix K such that K  $K^{T} = I$ , and therefore many solutions for matrix B exist, all valid since the elements of B have no physical significance as far as synthetic hydrology is concerned (Matalas, 1967).

Several techniques have been proposed for the solution of equation (B.12). Fiering (1964) and Matalas (1967) suggested the use of principal component analysis and Moran (1970) used canonical correlation analysis. Young (1968) assumed that B is a lower triangular matrix, based on the fact that  $C = B B^T$  is a symmetric matrix, and gave a unique recursive solution for the elements of B. Let us examine this case closely: (1)  $C = B B^{T}$  is symmetric for any B. The (i,j)th element of matrix C is

$$c_{ij} = \sum_{k=1}^{m} b_{ik} b'_{kj}$$
(B.16)

and the across the diagonal element is

$$c_{ji} = \sum_{k=1}^{m} b_{jk} b'_{ki}$$
(B.17)

where the prime denotes a transposed element. Thus,  $b'_{kj} = b_{jk}$  and  $b'_{ki} = b_{ik}$ , which implies that  $c_{ij} = c_{ji}$  and therefore C is symmetric for any B.

(2) That C is symmetric implies that m(m+1)/2 equations are required to specify it, and so m(m+1)/2 non zero elements of matric B are needed. Thus, since the (mxm) matrix B has  $m^2$  elements there are m(m-1)/2 elements that can be set to zero. So the assumption of a lower triangular matrix B is valid.

(3) The assumption of a lower triangular matrix B allows a recursive solution for the coefficients of B. This will be illustrated in the (2x2) case, and the reader is referenced to Young and Pisano (1968) for the general case.

$$\begin{bmatrix} b_{11} & 0 \\ & & \\ b_{21} & b_{22} \end{bmatrix} \begin{bmatrix} b_{11} & b_{21} \\ & & \\ 0 & b_{22} \end{bmatrix} = \begin{bmatrix} c_{11} & c_{12} \\ & & \\ c_{21} & c_{22} \end{bmatrix}$$

or

(B.18)

$$\begin{bmatrix} b_{11}^2 & b_{11} & b_{21} \\ & & & \\ b_{21} & b_{11} & (b_{21}^2 - b_{22}^2) \end{bmatrix} = \begin{bmatrix} c_{11} & c_{12} \\ & & \\ c_{21} & c_{22} \end{bmatrix}$$

from which

$$b_{11} = c_{11}$$

$$b_{21} = c_{21}/b_{11}$$

$$b_{22} = \sqrt{c_{11} - b_{21}^{2}},$$
(B.19)

with the constraints

$$c_{11} \ge 0$$
 and  $c_{22} \ge c_{12}^2 / c_{11}$  (B.20)
#### APPENDIX C DATA USED AND STATISTICS

Table C.l. 55 years of monthly rainfall data for the South Florida Station 6038.

### \*\*\*\*\* STATION 6038, MODRE HAVEN LOCK 1 \*\*\*\*\*

### Table C.2. 55 years of monthly rainfall data for the South Florida Station 6013.

\*\*\*\*\* STATION 6013, AVON PARK \*\*\*\*\*

1199933323456789012345678901234567890123456789012345678901 11999999999999999999999999999999999
001490197249934299181054299929000191202934395025785589922900005196927 0014904924864589042091920102927724199259235785589922900055969545 001490011042119344111140000191202810216257855899229000051965245
7407645055533062246031218867186640004233875110327763462246 8101313813813142607400185065659091565251508980545289563141 111420021851133401023000024211053336316446211220345289563141
29159797585011154598904008857444904747552082157199574608449403 291557975853111522541014000105611114575573100888752889754608449483 291497595978585345598904018944490485755755735008885751997344008449483
19485588042974424974944050055575112373400400188188 974494400394497441 194855880429744457440500055575112373400400188188 974444003541841 19465588049744140414401510446931007411463397800000000000000000000000000000000000
035745379535043375374640385034149914785148814448003604545795376030977651 0357453794518544830440334434814699147012015181374715451 035745319448104050349375074448146991470120151814448003604556000977651
8681394974770419101499747948888657977766088980921100445659796 9943727988069495779907970883199197105800898092100445659796 11114689057490105 111794060147898857488886579777666088980921100445659796 1117940798806945579007970889197105880865797776699161468905740057
9119588300993246667888092435754647473427938985459856569823645 3954914715881148587114908635754647406094462713956569823645 5354843411588711490863575464749793898545985459823645 115871490863575464247947938985455985459823645
9650750009977450192047097197400165910777298584288997192083259 965075621197649479506209744097698999900742858428899982259 1762119764999228801104709719774098999900742858428899982259 17621197649992288011920470971974098999900742858428899 17621197649992288011920470971974098999919208924599 17621197649992288011920470971974098984288999884288999919208832859 176211976499922880119209997444466
849244617597127488788789739108392109709088070175550447497296423380 3011874131962694533884687352688674242707777475607396422338 11117717774756073908844687352688674242707777475607396422338
80502804197774382355447906402685566135162641518363822894191 807449599198446661349799759888647472012446208992194320641116 9222309992198446661349799756612712192100420097221406211140
036588735740009 W496999W00404040013888879544750099199988 40501449900455140009 W4969730446509770540711376543445503919974040 01010030115000000100000110117100101000050001000010040000140101
10N41001N00041000N1110101010100N01010101

Note: -l indicates missing value.

# Table C.3. 55 years on monthly rainfall data for the South Florida Station 6093.

\*\*\*\*\* STATION 6093, FORT MYERS WSD AP \*\*\*\*\*

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## Table C.4. 55 years of monthly rainfall data for the South Florida Station 6042.

\*\*\*\*\* STATION 6042, CANAL PDINT USDA \*\*\*\*\*

Note: -1 indicates missing value.

## Table C.5. Monthly statistics of stations 6038, 6013, 6093, 6042.

***** STATION 6038 *****							
VARIABLE	N	MEAN	STANDARD DEVIATION	SKEWNESS	c. v.		
JAN FEB APR MAR JUN JUN JUC SEP DC DC DC DEC	55555555555555555555555555555555555555	1.927 1.8597 2.5597 4.5775 7.03577 7.0337 7.77,0379 1.379 1.457	3.063 1.368 2.456 1.818 2.584 3.358 2.897 4.085 3.073 1.283 1.555	5.014 0.664 1.762 0.674 1.032 0.646 1.008 1.008 1.081 1.532 1.975	159.002 72.854 94.854 56.482 49.644 40.420 41.193 53.983 53.983 53.042 106.686		
		**** STAT	ION 6013 ***	**			
VARIABLE	N	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.		
JAN FER APR JUL AUC SECT NDC DEC	000400055544 55555555555555555555555555	2.073 7187 22.7982 79882 79822 79822 79822 79822 7982 798	1.780 1.828 2.029 2.455 3.4694 3.148 3.732 2.488 1.551 1.135	1.344 0.798 0.864 1.073 1.129 0.793 1.487 0.716 0.798 1.833 0.727	85.0439 67.259 67.259 63.359 42.359 42.359 44.108 42.359 44.108 47.08 47.08 200 71.082 98.280 67.280		
		**** STAT	ION 6093 ***	**			
VARIABLE	N	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.		
JAN FEB AARR MADR JUL AUEP SCCV DEC	555555555555555555555555555555555555555	1.634 2.639 1.975 4.045 8.672 8.3553 8.3553 3.474 1.399	1.587 1.450 3.450 1.953 2.414 4.082 2.976 3.498 2.877 1.020	1.531 0.588 2.779 1.474 0.381 0.777 0.123 0.974 0.407 1.295 0.881 1.555	97.018 71.102 122.418 97.916 59.6155 34.3137 41.597 46.624 89.185 90.105		
		**** STAT	ICN 6042 ***	**			
VARIABLE	N	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.		
JAN FEB MAR APR JUL JUL SEP OCT NDC DEC	0004454555444 5555555555555555555555555	1.686 1.962 2.850 4.622 8.141 7.861 7.1802 4.702 4.702 1.818	1.488 1.475 2.2578 2.259 2.4108 2.4557 3.4552 3.1263 3.489 1.863	1.812 0.881 2.388 0.758 0.530 0.256 0.666 0.656 1.056 5.743 1.362	100.077 75.242 95.010 77.482 50.4446 43.353 39.4553 52.880 47.1632 176.912 102.457		

Table C.6. Station 6038--monthly statistics of the incomplete and estimated series--2% missing values.

	**** ST	ATION 603	38 ( 2%	MIS.	) ****	
VARIABLE	N	MEAN	STANDA DEVIAT	RD ION	C. V.	SKEWNESS
JEARRY JULGPTVC	55555555555555555555555555555555555555	1.885 1.96227 4.96377 4.5210 7.530307 7.530307 3.747 1.327 1.327	3.1.2.1.2.5.2.4.3.1.1.	034832985385 8839385385	163.191 71.646 93.646 72.498 55.855 56.762 46.513 41.798 53.983 82.017 92.781 108.781	5.083 0.643 1.744 1.0147 0.7147 0.753 0.753 0.753 1.081 1.138 1.385 1.943
	***** USING	THE MEAN	VALUE	( 2%	MIS. ) *****	
VARIABLE	MEAN	STAI DEV	NDARD IATION		C.V.	SKEWNESS
JAB FER APRY JULGP SEC NDC NDC	1.885 1.904 2.632 2.507 4.624 7.510 7.308 7.030 7.030 7.547 3.747 1.324 1.457		3. 048 1. 353 2. 441 1. 818 2. 559 3. 741 3. 336 2. 883 4. 085 3. 073 1. 217 1. 555		161. 666 70. 976 72. 746 72. 498 55. 336 49. 813 45. 643 41. 016 53. 983 82. 017 91. 911 106. 738	5.127 0.6499 0.6794 1.0754 0.754 0.7309 0.7309 0.781 1.1388 1.6597 1.977
***	** RECIPROCAL	DISTANCE	ES METH	00 (	2% MIS. ) **	***
VARIABLE	MEAN	STA! DEV	NDARD IATION		C. V.	SKEWNESS
JAB RABR RAPAY JULGP SECDO DE NDE	1.921 1.878 2.598 2.597 4.563 7.582 6.997 7.567 3.747 1.376 1.460		3.059 1.368 2.453 1.818 3.805 3.805 3.805 3.805 3.891 4.085 3.0738 1.556		159.245 72.821 94.413 72.498 50.929 50.883 41.310 53.983 82.017 92.825 106.582	5 039 0 464 1 674 1 0 671 0 672 0 764 1 087 0 764 1 138 1 527 1 968
	***** NORMAL	RATIO ME	ETHOD (	2% M	1IS. ) <del>****</del> *	
VARIABLE	MEAN	STAN DEV	NDARD		C. V.	SKEWNESS
JAB FEAR MAPAY JUL GP JUL SCCV ND C DD C	1.927 1.876 2.598 2.597 4.560 7.279 6.977 7.567 3.747 1.349 1.448		3.064 1.370 2.453 1.818 2.603 3.757 2.896 4.085 3.073 1.231 1.556		158.978 73.020 94.437 72.498 57.080 49.839 45.877 41.516 53.983 82.017 91.259 107.430	5:014 0:660 1:766 0:674 1:003 0:773 0:773 0:773 1:081 1:138 1:566 1:972
***	*** MODIFIED	WEIGHTED	AVERAG	E ( 2	2% MIS. ) ***	**
VARIABLE	MEAN	STAP DEV	NDARD		C. V.	SKEWNESS
JAN FEB APR MAPR JUN JUL SEP DCT NDC DEC	1.953 1.871 2.597 2.597 4.551 7.235 6.573 7.235 6.567 3.747 1.349		3.089 1.377 2.459 1.818 2.615 3.812 3.362 2.909 4.085 3.073 1.230 1.556		158. 165 73. 587 94. 897 72. 498 55. 339 46. 473 41. 779 53. 983 82. 017 91. 241 107. 386	4. 889 0. 643 1. 757 0. 674 0. 977 0. 684 1. 000 0. 773 1. 081 1. 138 1. 571 1. 986

Table C.7. Station 6038--monthly statistics of the incomplete and estimated series--5% missing values.

	** <b>*</b>	ATION 6038 ( 5	% MIS. ) *****	
VARIABLE	N	MEAN STAND DEVIA	ARD C.V. TION	SKEWNESS
JEARRYDUUGPTVC	5478 55451 557555555555555555555555555555555	1.885       3         1.725       1         2.358       1         4.521       2         7.411       3         7.355       3         7.474       4         3.790       3         1.344       1         1.457       1	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	1       5.083         1       0.616         0       1.804         3       0.664         0       0.853         7       0.719         6       0.784         3       0.676         7       1.147         1       1.31         7       1.643         5       1.975
	***** USING	THE MEAN VALUE	( 5% MIS. ) ***	<b>!</b> ₩.₩
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS
JAN FEB MAR APR JUL JUL AUG SECT NDV DEC	1.885 1.925 2.368 4.521 7.417 7.356 7.062 7.062 7.474 3.790 1.344 1.457	3.048 943 2.3457 1.645 2.5256 2.5293 2.8955 3.8955 3.0203 1.255	161.666 69.758 89.985 70.330 51.866 48.299 44.772 40.919 54.255 79.711 89.555 106.686	5.127 0.633 1.922 0.689 0.884 0.738 1.001 0.702 1.167 1.161 1.688 1.975
***	** RECIPROCAL	DISTANCES METH	HOD ( 5% MIS. )	****
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS
JAN FEB MAPR JUL AUEP JUL AUEP NDEC NDEC	1. 921 1. 867 2. 580 2. 429 4. 417 7. 613 7. 259 7. 039 7. 039 3. 837 1. 375 1. 457	3.059 1.370 2.426 1.786 2.414 3.772 2.899 4.299 4.290 1.270 1.555	159.265 73.3368 74.058 73.539 54.646 49.603 45.904 41.128 55.526 92.366 106.686	5.039 0.6832 0.667 0.6853 0.653 1.034 0.722 1.028 1.098 1.578 1.975
	***** NORMAL	RATIO METHOD	( 5% MIS. ) ****	÷¥
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS
JAN FEB MAR MAPR JUL JUL G SECT NDC NDC	1.927 1.856 2557 2.403 4.434 7.536 7.062 7.062 7.0691 3.773 1.335 1.457	3.064 1.377 2.423 1.752 2.398 3.3690 3.3695 2.890 4.365 2.8921 3.041 1.231 1.555	158, 978 74, 160 94, 741 72, 939 54, 081 48, 961 46, 590 40, 919 54, 883 80, 588 92, 198 106, 686	5.014 0.683 1.847 0.660 0.878 0.643 1.012 0.702 1.008 1.157 1.608 1.975
**1	*** MODIFIED	VEIGHTED AVERAG	E ( 5% MIS. ) *	***
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS
JAN FEB MAR APR JUN JUL SEP DCT NDC NDCC	1.953 1.849 2.405 4.403 7.584 7.197 7.019 3.816 1.349 1.457	3.089 1.387 2.459 1.773 2.4387 3.4387 3.4387 3.4387 3.4387 3.4387 3.4387 3.4387 3.4387 3.0925 3.0921 1.255	158.165 74.995 96.012 73.744 55.374 49.934 47.191 41.409 56.753 81.032 90.547 106.686	4.889 0.657 1.756 0.644 0.846 0.973 0.725 1.085 1.106 1.624 1.975

Table C.8. Station 6038--monthly statistics of the incomplete and estimated series--10% missing value.

	***++ STA	TION 603	8 ( 10% MIS	) <del>***</del> **	
VAR IABLE	N	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS
JAB FEAR APAYNL JUUPTJUC AUECTUC	50 472 549 481 550 491 550 551	1.848 1.924 2.505 4.4807 7.1820 7.1582 7.5826 1.318 1.486 1.318 1.486	3.108 1.374 2.463 2.468 3.748 3.748 3.748 2.748 2.748 2.740 4.124 4.976 1.176 1.575	168.167 71.410 78.180 73.066 54.985 47.931 45.803 41.064 54.393 80.318 89.225 107.372	5.328 0.705 1.051 0.631 1.055 1.067 0.677 1.078 1.261 1.498 1.943
	***** USING T	HE MEAN	VALUE ( 10%	MIS. ) *****	
VARIABLE	MEAN	STAI DEV	NDARD IATION	C. V.	SKEWNESS
JERRYNLGPTY AEAPAJJJAECD	1.848 1.923 2.569 2.564 4.488 7.807 7.223 7.123 7.1582 3.706 1.314		2.961 1.268 2.394 1.767 2.3001 3.451 2.301 3.151 2.3.969 2.835 2.835 1.117	160. 178 65. 926 95. 411 68. 873 51. 295 46. 120 43. 632 38. 632 38. 631 52. 341 76. 501 85. 021	5.572 0.760 1.765 0.665 1.120 0.700 1.118 0.736 1.117 1.318 1.568
***	** RECIPROCAL	DISTANCE	ES METHOD (	10% MIS. ) **	***
VAR IABLE	MEAN	STAN	NDARD IATION	C. V.	SKEWNESS
JABBRRY FEARRY JJJJGPTVC SCONDEC	1.872 1.876 2.608 4.430 7.621 7.435 7.158 7.581 3.746 1.323 1.445		3.025 1.366 2.462 1.941 2.468 3.808 3.313 2.832 3.032 4.042 3.032 1.156 1.553	161.608 72.829 94.405 55.572 49.967 44.566 39.562 52.631 80.926 87.370 107.435	5.225 0.7345 0.7452 0.9799 0.4861 0.715 1.020 1.421 2.007
	**** NORMAL	RATIO M	ETHOD ( 10%	MIS. ) *****	
VARIABLE	MEAN	STAN	NDARD IATION	C. V.	SKEWNESS
JABBRRY JFEARRY JJULGPT JJULGPT JJULGPT JOEC NDE	1.883 1.817 2.556 4.478 7.632 7.263 7.121 7.624 3.660 1.347 1.451		3.025 1.341 2.448 1.870 2.432 3.814 3.188 2.800 4.019 2.963 1.1753	160.648 73.800 94.535 73.141 54.064 49.972 43.889 39.326 52.721 80.954 87.209 107.011	5.217 0.803 1.775 0.478 0.9647 1.0446 0.753 1.044 1.192 1.382 2.000
**	*** MODIFIED V	NEIGHTED	AVERAGE (	10% MIS. ) ***	**
VARIABLE	MEAN	STAN DEVI	NDARD LAT LON	C. V.	SKEWNESS
JAN FEB MAR JUL JUL SEP DCD NDC	1 908 1.341 2.616 2.596 4.425 7.537 7.399 7.115 7.681 3.693 1.299 1.419		3. 145 1. 402 2. 478 1. 959 2. 532 3. 935 3. 359 3. 359 4. 084 3. 119 1. 129 1. 568	165.872 76.132 75.448 57.217 52.237 45.398 40.689 53.193 84.458 86.871 110.497	4. 645 0. 696 1. 728 0. 780 0. 947 0. 826 0. 783 1. 003 1. 155 1. 557 1. 971

Table C.9. Station 6038--monthly statistics of the incomplete and estimated series--15% missing values.

**<** STATION 6038 ( 15% MIS. ) *****						
VARIABLE	N MI	EAN STANDARD DEVIATION	c. v.	SKEWINESS		
JAN FEBR APR JUL SEP DCT DEC DEC	4653 477 478 505 477 496 49	1.927       3.284         1.750       1.449         2.448       2.099         2.448       1.857         4.674       2.490         7.727       3.854         7.111       3.467         7.245       2.954         7.245       3.954         3.456       3.003         1.488       1.343         1.488       1.378	170, 428 82, 791 85, 738 74, 749 53, 272 49, 876 48, 762 40, 777 51, 632 82, 138 90, 292 108, 686	4.910 0.875 1.171 0.749 0.887 0.592 1.101 0.664 0.673 1.182 1.417 1.975		
	***** USING THE	E MEAN VALUE ( 15	% MIS. ) ++++			
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS		
JAB AAB AAARYNLGPTVC SCCDC	1. 927 1. 750 2. 449 2. 484 4. 673 7. 728 7. 111 7. 244 7. 293 3. 656 1. 488 1. 470	2,998 1,308 1,851 1,714 2,298 3,303 2,667 3,475 2,831 1,226 1,506	155.539 74.733 75.631 69.007 49.174 46.450 36.813 47.657 77.431 82.403 102.471	5.338 0.984 1.314 0.808 0.956 0.631 1.152 0.730 0.725 1.247 1.540 2.085		
	** RECIPROCAL D	ISTANCES METHOD	( 15% MIS. ) *	****		
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS		
JAN FER MARRY JUL AUL JUGP TUL SECT NDE NDE	2.015 1.898 2.672 2.659 4.711 7.676 7.159 7.174 7.401 3.893 1.402 1.463	3.019 1.443 2.557 1.993 2.574 3.338 2.745 3.338 2.749 3.118 1.272 1.537	149.847 76.039 95.699 74.944 54.652 49.050 46.633 38.261 50.660 80.103 90.778 105.049	5. 144 0. 757 2. 011 0. 775 0. 775 0. 579 0. 734 0. 714 1. 048 1. 534 1. 979		
	***** NORMAL R	ATIO METHOD ( 157	( MIS. ) *****			
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS		
JAN FEB MARR MARR JUL AUGP SCCV NDC	2.017 1.808 2.621 2.571 4.750 7.585 7.129 7.130 7.130 7.130 7.412 1.463	3.029 1.370 2.377 1.878 2.599 3.362 2.739 3.362 2.739 3.718 2.989 1.278 1.538	150. 183 75. 774 90. 718 73. 030 54. 711 49. 208 47. 161 38. 419 50. 843 78. 337 90. 467 105. 085	5. 095 0. 812 1. 575 0. 425 0. 794 0. 429 1. 084 0. 757 0. 499 1. 009 1. 507 1. 994		
***	** MODIFIED WE	IGHTED AVERAGE (	15% MIS. ) ***	**		
VARIABLE	MEAN	STANDARD DEVIATION	C. V.	SKEWNESS		
JAN FEB MAR APR JUL SEP OCT NDC DEC	2.152 1.853 2.490 4.760 7.557 7.053 7.074 7.315 3.865 1.413 1.431	3.093 1.467 2.547 1.969 2.707 3.871 3.369 2.8835 3.857 3.857 3.203 1.248 1.553	143.705 79.169 97.056 51.220 47.772 39.965 52.859 82.859 88.354 108.479	4.701 0.844 1.981 0.564 0.797 0.529 1.113 0.700 0.766 1.120 1.623 1.956		

Table C.10. Station 6038--monthly statistics of the incomplete and estimated series--20% missing values.

	***** STA	TION 603	8 ( 20% MI	S. ) *****	
VARIABLE	N	MEAN	STANDARD DEVIATION	<b>c</b> . <b>v</b> .	SKEWNESS
JAN FEB AARR MARR JULG SECT NDC DEC	4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4 4	1.854 1.909 2.5797 4.797 7.3023 7.3023 3.841 1.3573	3. 240 1. 436 2. 112 2. 769 3. 720 3. 720 2. 824 4. 142 3. 210 1. 317 1. 703	174, 615 75, 213 87, 2541 72, 393 57, 730 53, 376 50, 916 40, 233 55, 024 83, 576 96, 582 108, 218	5.048 0.656 1.144 0.701 0.912 0.887 0.383 1.291 1.180 1.603 1.765
	**** USING T	HE MEAN V	VALUE ( 201	% MIS. ) 🕶 🛪 🛪 🛪	
VARIABLE	MEAN	STAN DEVI	NDARD IATION	C. V.	SKEWNESS
JABRRYNLGPTVC	1.856 1.909 2.411 2.572 4.797 7.307 7.307 7.307 7.307 7.528 3.841 1.363 1.573		2.991 1.296 1.906 1.738 2.442 3.439 3.319 2.462 3.653 2.878 1.148 1.501	161.109 67.886 79.547 50.905 47.067 45.430 35.064 48.524 75.446 84.213 75.482	5.434 0.720 1.2578 1.0237 0.9133 0.9133 0.94358 1.4358 1.820 1.820 1.981
**	*** RECIPROCAL	DISTANCE	S METHOD	( 20% MIS. ) +	****
VARIABLE	MEAN	STAN	NDARD IATION	C. V.	SKEWNESS
JABRRYN FEARRYNLGPTVC Abbrrynlgptvc	1.930 1.870 2.638 2.520 4.651 7.529 7.643 7.060 7.839 7.839 7.840 1.710		3.032 1.343 2.532 1.868 2.636 3.627 3.501 2.531 4.065 3.065 2.417 1.594	157.096 71.811 95.978 74.132 56.673 48.200 45.811 35.852 51.860 79.805 141.338 103.496	5 151 0 486 2 125 0 465 0 469 0 469 0 469 0 469 0 469 0 469 0 550 1 128 1 811
	***** NORMAL	RATIO ME	THOD ( 20%	4 MIS. ) *****	
VARIABLE	MEAN	STAN DEVI	IDARD ATION	<b>c</b> . <b>v</b> .	SKEWNESS
JAN FEAF MAPRY JUUL GPT JUUGPT JUUGPT VC SCCV CNDC	1.952 1.828 2.580 4.671 7.383 6.950 7.684 3.779 1.459 1.459		3.037 1.340 2.344 1.870 2.658 3.591 3.404 2.570 3.404 2.570 3.930 3.001 1.418 1.543	155.594 73.293 90.853 73.914 56.899 50.001 46.102 36.975 51.143 79.420 97.228 102.983	5. 110 0. 780 1. 645 0. 645 0. 859 0. 873 0. 428 1. 087 1. 232 1. 850 1. 748
**	*** MODIFIED W	EIGHTED	AVERAGE (	20% MIS. ) ***	***
VARIABLE	MEAN	STAN DEVI	DARD AT ION	C. V.	SKEWNESS
JAN FEB MAR MAR JUL JULG SEP DCV DEC DEC	2 027 1.823 2.591 4.705 7.283 7.572 6.941 7.829 3.738 1.510 1.494		3. 143 1. 371 2. 525 1. 886 2. 770 3. 670 3. 610 3. 610 4. 323 4. 323 3. 188 1. 655 1. 648	155.040 75.175 97.451 75.377 58.864 50.666 47.677 37.356 55.215 85.288 109.608 110.308	4.605 0.674 2.107 0.644 0.732 0.732 0.732 0.517 0.514 1.043 2.923 1.751

Table C.11. Lag-zero covariance matrices of the monthly rainfall series of stations 6013, 6093, 6042. All matrices are symmetric.

JAN:	3.169 2.393 2.343	2.518 1.723	2.848	,	FEB:	3.342 1.813 1.496	2.101 1.505	2.175
MAR:	4.022 3.501 2.677	10.275 6.921	6.254	,	APR:	4.881 1.959 3.545	3.814 2.378	5.190
MAY:	7.047 3.610 3.539	5.826 2.871	7.071	,	JUN:	13.647 8.235 7.373	16.666 7.138	16.865
JUL:	13.425 3.655 3.378	8.854 2.546	11.615	,	AUG:	9.907 -0.628 1.630	12.177 0.260	8.138
SEP:	13.928 9.080 5.913	15.902 6.443	17.022	,	ОСт:	6.189 5.032 4.516	8.278 5.896	10.004
NOV:	2.406 0.822 1.395	1.098 0.547	12.174	,	DEC:	1.289 1.045 1.145	1.588 1.510	3.471

Table C.12. Normality transformations applied on the monthly rainfall data of Station 6038.

	***** STATION	6038 ( NO TRA	NSFORMATION ) +	****
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JABB JFEARRY JUULGP AMADA JUULGP AUST AUST AUST AUST AUST AUST AUST AUST	1.927 1.878 2.595 2.507 4.575 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.033 7.035 7.057 7.035 7.0577 7.0577 7.0577 7.0577 7.0577 7.0577 7.0577 7.0577 7.0577 7.0577 7.	3.043 1.348 2.584 3.584 3.358 2.897 4.085 3.073 1.283 1.555	5.014 0.664 1.762 0.674 1.032 0.646 1.084 1.084 1.081 1.138 1.535 1.975	159       002         72.       856         94.       631         72.       498         56.       482         49.       420         53.       983         53.       983         62.       017         93.       042         106.       686
*****	STATION 6038	( LDGARITHMIC	TRANSFORMATION	) <del>=====</del>
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JAN FEB APR MAY JJJL AUG SECT AUG SECT DO DEC	-0.009 0.103 0.188 0.218 0.573 0.822 0.809 0.810 0.814 0.385 -0.088 -0.068	0. 532 0. 466 0. 521 0. 251 0. 251 0. 227 0. 184 0. 227 0. 184 0. 253 0. 488 0. 519 0. 479	-0. 266 -1. 224 -0. 936 -1. 349 -0. 164 -0. 875 -1. 034 -0. 226 -0. 645 -1. 330 -0. 755 -0. 173	-5598.555 452.424 276.239 230.355 42.276 29.918 28.124 22.643 31.143 126.596 -586.139 -706.420
	***** STAT	IDN 6038 ( POWE	ER=0.25 ) *****	÷
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	c. v.
JAN FEB MAR MAR JUL JUL AUG SET NCV DEC	1.040 1.076 1.161 1.175 1.421 1.606 1.603 1.614 1.272 0.990 0.998	0. 314 0. 256 0. 312 0. 263 0. 203 0. 218 0. 199 0. 168 0. 199 0. 168 0. 227 0. 316 0. 268 0. 270	0.738 -0.540 -0.108 -0.678 0.143 -0.355 -0.454 -0.140 -0.334 -0.130 0.364	30. 141 23. 382 26. 870 24. 109 14. 281 13. 449 12. 387 10. 457 14. 051 24. 436 27. 100 27. 061
	***** STAT	ION 6038 ( POWE	ER=0.35 ) ****	÷
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JAN FEB MAPR MAPR JUL AUEP JUL AUEP NDC NDC DDC DDC	1.083 1.154 1.274 1.274 1.644 1.975 1.949 1.949 1.942 1.965 1.456 1.006 1.017	0.461 0.361 0.407 0.328 0.328 0.366 0.332 0.283 0.382 0.480 0.382 0.480 0.370 0.383	1.219 -0.327 0.163 -0.443 0.264 -0.180 -0.238 0.121 0.040 -0.060 0.110 0.581	42 593 31 454 31.454 31.909 19.933 18.530 17.051 14.592 19.458 32.927 36.724 37.667
	***** STAT	IDN 6038 ( SQUA	RE ROOT ) ****	÷¥
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JAN FEAR MAPRY JJJJLGP AUSCOP AUSCONDEC	1. 178 1. 265 1. 442 1. 459 2. 059 2. 671 2. 618 2. 578 2. 578 1. 768 1. 051 1. 067	0.740 0.533 0.724 0.620 0.584 0.673 0.625 0.537 0.727 0.725 0.529 0.529	2.043 -0.048 0.540 -0.117 0.445 0.053 0.265 0.282 0.285 0.285 0.285 0.282 0.461 0.908	62.830 42.115 50.192 42.466 28.373 25.944 23.873 20.760 27.446 44.953 50.341 53.473

### Table C.13. Statistics of the estimated series -- univariate model

	**** UNIVARI	ATE MODEL ( 10%	MIS. ) ****	
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	c. v.
JEAR FEARRY JUDU JUDU SCCVC	1.854 1.891 2.503 2.499 4.504 7.809 7.185 7.067 7.563 3.594 1.314 1.475	2.980 1.278 2.402 1.805 2.305 3.601 3.161 2.809 3.975 2.871 1.132 1.539	5.462 0.815 1.743 0.701 1.095 0.697 1.141 0.780 1.127 1.369 1.515 2.019	160.709 67.608 95.956 72.242 51.181 46.120 43.995 39.757 52.555 79.865 86.162 104.304
	**** UNIVARI	ATE MODEL ( 20%	MIS. ) ****	
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JAB FEAR APRY JUUL GP AUNL AUEP CC T	1.777 1.846 2.523 4.713 7.199 7.216 6.961 7.420 7.719	2.997 1.303 1.914 1.743 2.449 3.449 3.325 2.465 3.455 3.459 2.910	5.480 0.854 1.364 0.828 1.119 1.009 1.062 0.510 1.531 1.408	168.623 70.585 81.959 69.053 51.959 47.865 46.072 35.406 49.316 78.235

## Table C.14. Statistics of the estimated series -- bivariate model

	*** BIVARIA	TE MODEL ( 10%	MIS. ) ****	
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	С. У.
JABRRY JEBRRY JULGPTVC JOUECTVC SCOZD	1.825 1.867 2.483 2.534 4.491 7.749 7.220 7.110 7.523 3.592 1.293 1.459	2.972 1.287 2.398 1.781 2.327 3.612 3.166 2.784 3.979 2.863 1.123 1.539 ATE MODEL ( 20%	5.532 0.841 1.976 0.658 1.086 0.740 1.104 0.779 1.154 1.392 1.392 1.599 2.052 MIS. ) ****	162 862 68 866 96 605 51 825 46 811 43 859 39 159 52 809 86 839 105 450
VARIABLE	MEAN	STANDARD DEVIATION	SKEWNESS	C. V.
JABRRRYN JEBRRYYN JUULGPTYC Secou Decou Decou	1.798 1.835 2.353 2.551 4.759 7.280 7.263 6.941 3.786 1.351	3.013 1.337 1.775 2.490 3.490 3.353 2.353 2.353 2.729 2.941 1.159 1.527	5.374 0.790 1.283 0.763 1.005 0.897 0.994 0.596 1.404 1.294 1.294 1.926	167 566 728 847 828 5527 459 55332 46 2537 46 239 46 239 46 239 77 684 87 456

#### APPENDIX D

#### COMPUTER PROGRAMS

#### RAEMV-U

(Recursive Algorithm for the Estimation of Missing Values - Univariate Model)

#### Input

The program inputs the time series; the parameters of the normality transformation to be performed (power transformation); the number of gaps (not necessarily the number of missing values unless all the gaps are singles); and for each gap the starting and ending point (counting starts from the first value in the series). For the first iteration the missing values in the original series (usually indicated by a code or by a negative value) are initialized to zeroes or to some other desired initial estimates.

#### Program Description

The main program reads the input data and then subsequently calls subroutine ARMA (each call corresponds to one iteration). Subroutine ARMA performs the following calculations each time it is called:

- The input series is transformed to normal (using the selected transformation) and stationary (by subtracting the monthly means and dividing by the standard deviations).
- (2) The mean, variance, autocovariance function (ACVF), autocorrelation function (AGF) and partial autocorrelation function (PACF) of the transformed series are computed by calling the IMSL subroutine FTAUTO.
- (3) Preliminary estimates of the p AR parameters, and q MA parameters are computed by calling the IMSL subroutines FTARPS and FTMPS subsequently.
- (4) Maximum likelihood estimates (MLE) of the AR and MA parameters are computed and the residual series is calculated by calling the IMSL subroutine FTMXL.
- (5) The mean, variance, ACVF, ACF and PACF of the residual series are computed by calling the IMSL subroutine FTAUTO.
- (6) The parameters of the fitted model (MLE) are used to estimate the missing values in all the gaps by the Box-Jenkins minimum mean square error forecasting procedure.
- (7) The inverse normality and stationarity transformations are performed on the series and the estimated complete series is output.

The estimated series (output from the first call) now becomes the input series for the second call and the above seven steps are repeated. The subroutine ARMA is called as many times as needed until stabilization of the parameter estimates and of the missing values estimates occur. The program is initialized to five calls (more can be easily added as needed), and a stabilization check for the parameters is provided so that the iterations stop when the two parameters remain constant to the second decimal place.

The computation and printing of the ACVF, ACF and PACF of the transformed and residual series (steps 2 and 5) are

not necessary and can be eliminated from the program without any problem. However, their inclusion permits the checking of the goodness of the fitted model at each iteration by diagnostic checking applied on the residuals. A listing of the program in FORTRAN follows.

#### RAEMV-B

(Recursive Algorithm for the Estimation of Missing Values - Bivariate Model)

The special case of having only the one series incomplete and the other complete will be considered here. However, the program can be easily modified to include the case of having both the series incomplete.

#### Input

The program inputs the two time series, the parameters of the normality transformation to be performed on each series, the number of gaps and the position of each gap for the incomplete series. The missing values in the incomplete series are initialized to zeros or to some other values.

#### Program description

The main program reads the input data and then subsequently calls subroutine BIVAR (each call corresponds to one iteration). Subroutine BIVAR performs the following each time is called:

- (1) The two input series are transformed to normal and stationary by calling subroutine STAT.
- (2) The lag-zero and lag-one autocovariances and cross-covariances of the two series are computed by calling the IMSL subroutine FTCRXY.
- (3) The parameter matrices A and B are calculated. Inversion and multiplication of matrices are performed by the IMSL subroutines LINV2F, VMULFF and VMULFP.
- (4) The parameter matrices A and B are used to estimate the missing values of the incomplete series.
- (5) The inverse normality and stationarity transformations are performed on the two series, and the estimated complete series is output.

The estimated series (output from the first call) now becomes the input series for the second call and the above five steps are repeated until stabilization of the matrices A and B occurs. No check for stabilization is provided by the program (eight values must be checked simultaneously) but instead the subroutine is called for a prefixed number of times. A listing of the computer program in FORTRAN follows.

ούουοοοούο \_\_\_\_\_ PROGRAM RAEMV-U RECURSIVE ALGORITHM FOR THE ESTIMATION OF MISSING VALUES - UNIVARIATE MODEL DIMENSION RAIN(60, 12), NYEAR(60) DIMENSION VRAIN(800), EIRAIN(60, 12), VRAIN1(800), E2RAIN(60, 12), VRAIN2(800), E3RAIN(60, 12), VRAIN3(800), E4RAIN(60, 12), VRAIN5(800), E6RAIN(60, 12), VRAIN5(800), E6RAIN(60, 12), VRAIN6(800) DIMENSION LI(200), LG(200), ISI(200), IEI(200) COMMON/A/ ID, NYEAR COMMON/B/ N, C, P COMMON/C/ NG, ISG(200), IEG(200) 000000000000 READ INPUT PARAMETERS HEADER. TITLE N.....NUMBER OF YEARS NG....NUMBER OF GAPS LI....LENGTH OF INTEREVENT LG....LENGTH OF GAP C.P....PARAMETERS OF THE TRANSFORMATION TRANSFORMED SERIES Y=(X+C)\*\*P READ(5,10) HEADER FORMAT(20A4) c<sup>10</sup> c<sup>20</sup> READ(5,20) C,P FURMAT(2F5,2) READ(5,30) N, NG FORMAT(2(14/)) с<sup>30</sup> DD 5 I=1,NG READ(5,40) LI(I),LG(I) FORMAT(2I4) 5 40 C READ(10, 110) (ID, (NYEAR(I), (RAIN(I, J), J=1, 12)), I=1, N) FORMAT(A4, I3, 1X, 12F6. 2) 110 FROM THE INPUT VARIABLES LI AND LG TWO ARRAYS OF LENGTH NG ARE COMPUTED. THEN THE STARTING POINT OF THE KTH GAP IS IBG(K) AND THE ENDING POINT IS IEG(K). 000000 ISI(1)=1 IEI(1)=LI(1) ISG(1)=IEI(1)+1 IEG(1)=ISG(1)+LG(1)-1 DO 11 I=2.NG ISI(I)=IEG(I-1)+1 IEI(I)=ISI(I)+1 IEG(I)=IEI(I)+1 IEG(I)=ISG(I)+LG(I)-1 CONTINUE c<sup>11</sup> 15 C C C WRITE(6,15) HEADER FORMAT(20A4,///) PRINT THE POSITION OF THE GAPS FOR CHECKING 100 C C C WRITE(6,100) (I,ISG(I),IEG(I),I=1,NG) FORMAT(316) INITIALIZE THE MISSING VALUES TO ZERO DD 50 I=1,N DD 50 J=1,12 IF(RAIN(I,J).EQ.-1) RAIN(I,J)=0. <sup>50</sup> CONTINUE

SUBROUTINE ARMA IS CALLED TO FIT AN ARMA(P,Q) MODEL TO THE INPUT SERIES. THE PARAMETERS OF THE MODEL ARE USED FOR THE ESTIMATION OF THE MISSING VALUES. CCCC CALL ARMA(RAIN, VRAIN, EIRAIN, VRAIN1, PHI1, THETA1) CALL ARMA(EIRAIN, VRAIN1, E2RAIN, VRAIN2, PHI2, THETA1) CALL ARMA(E2RAIN, VRAIN2, E3RAIN, VRAIN3, PHI3, THETA3) IF((PHI3-PHI2). LE. O. 001. AND. (THETA3-THETA2). LE. O. 001) GU 0979 CALL ARMA(E3RAIN, VRAIN3, E4RAIN, VRAIN4, PHI4, THETA4) IF((PHI4-PHI3). LE. O. 001. AND. (THETA4-THETA3). LE. O. 001) GU TO 979 CALL ARMA(E4RAIN, VRAIN4, E5RAIN, VRAIN5, PHI5, THETA5) IF((PHI5-PHI4). LE. O. 001. AND. (THETA5-THETA4). LE. O. 001) GU TO 979 CALL ARMA(E5RAIN, VRAIN5, E6RAIN, VRAIN6, PHI6, THETA6) C C 777 STOP ούοοοοοοο ----- SUBROUTINE ARMA ------SUBROUTINE ARMA FITTS AN ARMA(P,G) MODEL TO THE INPUT SERIES EACH TIME IS CALLED. THE MISSING VALUES ARE ESTIMATED BY THE BOX-JENKINS FORECASTING PROCEDURE AND THE ESTIMATED SERIES IS SAVED TO BE THE INPUT SERIES TO THE NEXT CALL. SUBROUTINE ARMA(RAIN, VRAIN, ERAIN, EVRAIN, PHI1, THETA1) С REAL MEAN(13), MP, LP DIMENSION ERAIN(60, 12), EVRAIN(800), Z(800) DIMENSION RAIN(60, 12), NYEAR(60), IND(8), PHI(11), THETA(11), SUMSG1(11, 11), ACV(300), AC(300), PACV(300), VRAIN(800), TEMP(800), WKAREA(1600), AC(300), GR(1600) DIMENSION YTOTAL(61), STD(13) COMMON/A/ ID,NYEAR COMMON/A/ N, C, P COMMON/C/ NG, ISG(200), IEG(200) CCCC INPUT THE VALUES OF PHI AND THETA FOR WHICH YOU WANT THE SUM OF SQUARES SURFACE TO BE CALCULATED. DATA PHI/- 5, - 4, - 3, - 2, - 1, 0, , 1, 2, 3, 4, 5/ DATA THETA/0, , 1, 2, 3, 4, 5, 6, 7, 8, 9, 1, / CCCC PRINT OUT THE ORIGINAL SERIES THIS IS THE SERIES THAT EACH ITERATION STARTS WITH. 112 2 111 C C \$ COMPUTE POWER TRANSFORMATION OF THE SERIES DO 10 I=1,N DO 10 J=1,12 RAIN(I,J)=(RAIN(I,J)+C)\*\*P CONTINUE 25 25 COMPUTE YEARLY TOTALS DO 22 I=1,N YTOTAL(I)=0. DO 23 J=1,12 YTOTAL(I)=YTOTAL(I)+RAIN(I,J) CONTINUE CONTINUE 232 C C C C C C COMPUTE MONTHLY MEANS AND STANDARD DEVIATIONS DD 32 J=1,12 MEAN(J)=0. STD(J)=0. DD 34 I=1,N MEAN(J)=MEAN(J)+RAIN(I,J)/FLOAT(N) STD(J)=STD(J)+RAIN(I,J)\*\*2 CONTINUE 34 32 c CONTINUE DO 36 I=1,12

```
STD(I)=((STD(I)-MEAN(I)**2*FLOAT(N))/(FLOAT(N)-1.))**0.5
CONTINUE
   с<sup>36</sup>
               MEAN(13)=0.

STD(13)=0.

DD 38 I=1,N

MEAN(13)=MEAN(13)+YTOTAL(I)/FLOAT(N)

STD(13)=STD(13)+YTOTAL(I)**2

CONTINUE

STD(13)=((STD(13)-MEAN(13)**2*FLOAT(N))/(FLOAT(N)-1.))**0.5
      38
    C
C
C
                NOW STANDARDIZE THE MONTHLY SERIES
                DD 42 I=1,N
DD 42 J=1,12
RAIN(I,J)=(RAIN(I,J)-MEAN(J))/STD(J)
CONTINUE
      42
    С
                STORE THE MATRIX SERIES IN A VECTOR SERIES
                DD 30 I=1,N
DD 30 J=1,12
K=J+(I-1)*12
VRAIN(K)=RAIN(I,J)
CONTINUE
     30
    C
C
C
C
C
C
                COMPUTE AC, PAC, AND ACV OF THE SERIES USING SUBROUTINE FTAUTO
               L=30
CALL FTAUTD(VRAIN, NN, L, L, 7, AMEAN, ACV(1), ACV(2), AC(2),
PACV(2), WKAREA)
    CCC
                SET AC AND PACY OF LAG ZERO TO ONE
   SSQ=0.

LP1=L+1

DD 40 I=1,LP1

IM1=I-1

WRITE(6,140) IM1,AC(I),PACV(I)

SSQ=SSQ+AC(I)**2

CONTINUE

SSQ=SSQ-1

FORMAT(3X, I5,2F15.7)

WRITE(6,145) SSQ

FORMAT(///,2X, SUM OF AC**2 (NOT INCLUDING LAG 0)',F15.6)
     40
145
C
C
C
C
C
               PRELIMINARY ESTIMATE OF AR PARAMETER AND OVERALL MA CONSTANT USING SUBROUTINE FTARPS
               CALL FTARPS(ACV,AMEAN,1,1,ARPS,PMAC,WKAREA)
WRITE(6,150) ARPS,PMAC
FORMAT(///,2X,'----SUBRDUTINE FTARPS-----',//,
1 5X,'ESTIMATE OF AR PARAMETER (ARPS).....',F15.6,//,
2 5X,'OVERALL MA CONSTANT (PMAC).....',F15.6,//)
     150
   CCCC
               ESTIMATE MA PARAMETER AND WHITE NOISE VARIANCE USING SUBROUTINE FIMPS
               CALL FTMPS(ACV, ARPS, 1, 1, PMAS, WNV, WKAREA, IER)
WRITE(6,155) PMAS, WNV
FORMAT(///,2x,'----SUBROUTINE FTMPS-----',//,
1 5X,'ESTIMATE OF MA PARAMETER (PMAS)....',F15.6,//,
2 5X,'WHITE NOISE VARIANCE (WNV).....',F15.6)
     155
             5
   CCC
               STORE VECTOR VRAIN BECAUSE FTMXL WILL DESTROY IT
               DD 45 I=1, NN
TEMP(I)=VRAIN(I)
CONTINUE
     45
   CCCC
               GENERATE SUM OF SQUARES SURFACE OF THE RESIDUALS
LET ITHETA AND IPHI BE THE ROW AND COLUMN NUMBER
CORRESPONDING TO THE MINIMUM SUM OF SQUARES OF
```

```
c
                    THE RESIDUALS
                   ITHETA=1

IPHI=1

DD 50 I1=1,11

DD 55 I2=1,11

ETA=TEMP(2)-PHI(I2)*TEMP(1)

SUMSQ1(I1,I2)=ETA**2

DD 60 I3=3,NN

ETA1=TEMP(I3)-PHI(I2)*TEMP(I3-1)+THETA(I1)*ETA

ETA=ETA1

SUMSQ1(I1,I2)=SUMSQ1(I1,I2)+ETA1**2

CDNTINUE

IF(SUMSQ1(I1,I2)=SUMSQ1(ITHETA,IPHI)) GD TD 55

ITHETA=I1

IPHI=I2

CONTINUE
      60
55
50
C
C
C
                   CONTINUE
                   WRITE OUT THE SUM OF SQUARES SURFACE OF THE RESIDUALS
                  WRITE(6,160)

FORMAT(1H1,//,50X,'TABLE 2',//,15X,'SUM DF SQUARES DF THE',

* 'RESIDUALS OF THE STANDARDIZED TRANSFORMED SERIES',

#//,52X,'PHI')

WRITE(6,165) (PHI(I), I=1,11)

FORMAT(5X,'THETA',2X,11(3X,F5.2,1X)/)

WRITE(6,170) (THETA(I),(SUMSQ1(I,J),J=1,11),I=1,11)

FORMAT((5X,F5.2,3X,11(FB.2,1X)))
      160
                 $
      165
     170
   000000
                   LET THE VALUES OF ARPS AND PMAS THAT LED TO MINIMUM SUM
OF SQUARES OF RESIDUALS BE INPUT TO SUBROUTINE FTMXL.
COMPUTE IMPROVED ESTIMATES OF ARPS, PMAC, PMAS AND WNV
USING SUBROUTINE FTMXL
                   DATA IND/0, 1, 1, 0, 75, 4, 1, 3/
IND(1)=NN
PMAS=THETA(ITHETA)
ARPS=PHI(IPHI)
   C
C
C
C
                    IF PHI=1 SET PHI=0. 99 SINCE PHI=1 IS NOT DESIRABLE
                   IF(PMAS.EG.1.) PMAS=0.99
WRITE(6,180) ARPS, PMAS
FORMAT(///,2%, THE SUM OF SQUARES OF RESIDUALS IS MINIMUM',
FOR: ', //,4%, 'ARPS= ',F15.6,/,4%, 'PMAS= ',F15.6,/)
      180
                   CALL FTMXL(TEMP, IND, ARPS, PMAS, PMAC, WNV, GR, A, IER)
WRITE(6, 190) ARPS, PMAS, PMAC, WNV
FORMAT(///, 2X, '----SUBROUIINE FTMXL----',//,
1 5X, 'ESTIMATE OF AR PARAMETER (ARPS).....', F15.6,//,
2 5X, 'ESTIMATE OF MA PARAMETER (PMAS).....', F15.6,//,
3 5X, 'OVERALL MA CONSTANT (PMAC).....', F15.6,//,
4 5X, 'WHITE NOISE VARIANCE (WNV)....', F15.6)
PH11=ARPS
THETA1=PMAS
   С
      190
                 123
                 ā
   CCCC
                   FIND AC AND PACY OF RESIDUALS THAT ARE STORED IN THE VECTOR (A) AS DUTPUT FROM SUBROUTINE FTMXL
                   CALL FTAUTO(A, NN, L, L, 7, AMEAN, ACV(1), ACV(2), AC(2),
PACV(2), WKAREA)
                 $
                   AC(1)=1.
PACV(1)=1.
   CCC
                    WRITE OUT AC AND PACY . COMPUTE SUM OF AC**2
                  SSGR=0.

WRITE(6,200)

FORMAT(1H1,///,15X, 'RESIDUAL SERIES ',///)

WRITE(6,135)

DD 70 I=1,LP1

IM1=I-1

WRITE(6,140) IM1,AC(I),PACV(I)

SSGR=SSGR+AC(I)**2

CONTINUE

SSGR=SSGR-1.

WRITE(6,145) SSGR
      200
      70
   С
                   DO 15 I=1, NN
EVRAIN(I)=VRAIN(I)
  c<sup>15</sup>
                   GENERATE RANDOM NUMBERS N(0,1) TO BE USED FOR THE FORECASTING
   c
```

```
С
                                     DSEED=123457.DO
CALL GGNML(DSEED, NN, Z)
       C
C
                                   DO 20 I=1,NG

I1=ISG(I)

I2=IEG(I)

K=I2-I1+1

IF(K.GT.1) ©O TO 51

EVRAIN(I1)=PHI1*VRAIN(I1-1)-THETA1*Z(I1-1)

GO TO 20

EVRAIN(I1)=PHI1*VRAIN(I1-1)-THETA1*Z(I1-1)

DO 31 L=2,K

EVRAIN(I1+L-1)=PHI1*EVRAIN(I1+L-2)

CONTINUE
             51
5.
31
20
C
C
C
C
                                      APPLY THE INVERSE TRANSFORMATIONS ON THE SERIES.
                                    PP=1/P

DD 61 I=1,N

K1=(I-1)*12+1

K2=I*12

DD 71 L=K1,K2

J=L-(I-1)*12

ERAIN(I,J)=(EVRAIN(L)*STD(J)+MEAN(J))**PP

CONTINUE

CONTINUE
       71
61
C
    61 CONTINUE

C

RETURN

END

//GO. SYSIN DD *

***** STATION 6038 - UNIVARIATE MODEL *****

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2 1

2 5

1 2

3 0

2 //GO. FT10F001 DD DSN=UF. B0063401. S7. C60381, DISP=(OLD, KEEP)
```

 $\sim$ 

с с	
C C C	PROGRAM RAEMV-B
с с с с	RECURSIVE ALGORITHM FOR THE ESTIMATION OF MISSING VALUES - BIVARIATE MODEL
	DIMENSION RAIN1(60,12), VR1(800), RAIN2(60,12), VR2(800), 1 E1R1(60,12), V1R1(800), E1R2(60,12), V1R2(800), A1(2,2), 2 B1(2,2), M01(2,2), M11(2,2) DIMENSION E2R1(60,12), V2R1(800), E2R2(60,12), V2R2(800), 1 A2(2,2), B2(2,2), M02(2,2), M12(2,2) DIMENSION E3R1(60,12), V3R1(800), E3R2(60,12), V3R2(800), 1 A3(2,2), B3(2,2), M03(2,2), M13(2,2) DIMENSION E4R1(60,12), V4R1(800), E4R2(60,12), V4R2(800), 1 A3(2,2), B4(2,2), M04(2,2), M14(2,2) DIMENSION E5R1(60,12), V3R1(800), E5R2(60,12), V4R2(800), 1 A4(2,2), B4(2,2), M04(2,2), M14(2,2) DIMENSION E5R1(60,12), V5R1(800), E5R2(60,12), V5R2(800), 1 A4(2,2), B5(2,2), M05(2,2), M15(2,2) DIMENSION LI(200), LG(200), ISI(200), IEI(200) COMMON/A/ ID1, ID2, NYEAR(60) COMMON/A/ N, C, P COMMON/C/ NG, ISG(200), IEG(200)
	READ INPUT PARAMETERS HEADER. TITLE NNUMBER OF YEARS NGNUMBER OF GAPS LILENGTH OF INTEREVENT LGLENGTH OF GAP C,PPARAMETERS OF THE TRANSFORMATION TRANSFORMED SERIES Y=(X+C)**P
, 10	READ(5,10) HEADER FORMAT(20A4)
20	READ(5,20) C,P FORMAT(2F5,2)
30	READ(5,30) N,NG FORMAT(2(14/))
5 40	DO 5 I=1,NG READ(5,40) LI(I),LG(I) FORMAT(2I4)
, 100	READ(10,100) (ID1, (NYEAR(I), (RAIN1(I,J), J=1,12)), I=1,N) READ(11,100) (ID2, (NYEAR(I), (RAIN2(I,J), J=1,12)), I=1,N) FORMAT(A4, I3, 1X, 12F6.2)
	FROM THE INPUT VARIABLES LI AND LO TWO ARRAYS DF LENGTH NG ARE COMPUTED. THEN THE STARTING POINT OF THE KTH GAP IS ISG(K) AND THE ENDING POINT IS IEG(K).
11	ISI(1)=1 IEI(1)=LI(1) IEG(1)=IEI(1)+1 IEG(1)=IEG(1)+LG(1)-1 DO 11 I=2,NG ISI(I)=IEG(I-1)+1 IEI(I)=IEI(I)+LI(I)-1 IEG(I)=IEG(I)+LG(I)-1 CONTINUE
15	WRITE(6,15) HEADER FORMAT(2044,///)
	PRINT THE POSITIONS OF THE GAPS FOR A CHECK
101	WRITE(6,101) (I,ISG(I),IEG(I),I=1,NG) FORMAT(316)
	INITIALIZE THE MISSING VALUES OF THE INCOMPLETE

С DO 60 I=1,N DO 60 J=1,12 IF(RAIN1(I,J).EQ.-1) RAIN1(I,J)=0. 60 C C C CONTINUE PRINT OUT THE SERIES WHICH IS TO BE ESTIMATED WRITE(6,66) WRITE(6,102) (ID1,(NYEAR(I),(RAIN1(I,J),J=1,12)),I=1,N) FORMAT(1X,A4,I3,1X,12F6.2) FORMAT(1H1) 102 66 000000 SUBROUTINE BIVAR IS CALLED TO FIT A BIVARIATE AR(1) MODEL TO THE TWO INPUT SERIES. IT ESTIMATES ALSO THE MISSING VALUES OF THE ONE SERIES AND SAVES IT TO BE INPUT TO THE NEXT CALL CALL BIVAR (RAIN1, VR1, RAIN2, VR2, E1R1, V1R1, E1R2, V1R2, A1, B1, M01, M11) CALL BIVAR (E1R1, V1R1, E1R2, V1R2, E2R1, V2R1, E2R2, V2R2, A2, B2, M02, M12) CALL BIVAR (E2R1, V2R1, E2R2, V2R2, E3R1, V3R1, E3R2, V3R2, A3, B3, M03, M13) CALL BIVAR (E3R1, V3R1, E3R2, V3R2, E4R1, V4R1, E4R2, V4R2, A4, B4, M04, M14) CALL BIVAR (E4R1, V4R1, E4R2, V4R2, E5R1, V5R1, E5R2, V5R2, A5, B5, M05, M15) С STOP END 0000000000000 ----- SUBROUTINE BIVAR -----SUBROUTINE BIVAR FITS A BIVARIATE AR(1) MODEL TO THE TWO INPUT SERIES EACH TIME IS CALLED. IT ESTIMATES ALSO THE MISSING VALUES OF THE ONE SERIES AND THE ESTIMATED SERIES IS SAVED TO BE INPUT TO THE NEXT CALL SUBROUTINE BIVAR (RAIN1, VR1, RAIN2, VR2, ERAIN1, EVR1, ERAIN2, EVR2, A, B, M0, M1) с DIMENSION WKAREA(200) DIMENSION RAIN1(60,12), VR1(800), RAIN2(60,12), VR2(800) DIMENSION ERAIN1(60,12), EVR1(800), ERAIN2(60,12), EVR2(800) DIMENSION XM1(12), XM2(12), STD1(12), STD2(12) DIMENSION A(2,2), B(2,2), C(2,2), D(2,2) REAL M0(2,2), M1(2,2), M0INV(2,2) COMMON/A/ ID1, ID2, NYEAR(60) COMMON/B/ N, C,P COMMON/C/ NG, ISG(200), IEG(200) COMMON/C/ XM1, XM2, STD1, STD2, X1, X2, ST1, ST2 0000 CALL SUBROUTINE STAT TO NORMALIZE AND STANDARDIZE THE SERIES AND COMPUTE THE STATISTICS CALL STAT(RAIN1, XM1, STD1, VR1, X1, ST1) CALL STAT(RAIN2, XM2, STD2, VR2, X2, ST2) 0000 CALL THE IMSL SUBROUTINE FICRXY TO COMPUTE AUTO-AND CROSS-COVARIANCES OF THE SERIES CALL FTCRXY(VR1, VR2, N, X1, X2, O, N, C12O, IER) CALL FTCRXY(VR1, VR1, N, X1, X1, -1, N, C111, IER) CALL FTCRXY(VR2, VR2, N, X2, X2, -1, N, C221, IER) CALL FTCRXY(VR2, VR2, N, X1, X2, -1, N, C121, IER) CALL FTCRXY(VR2, VR1, N, X2, X1, -1, N, C211, IER) С MO(1,1)=1. MO(2,2)=1. MO(1,2)=C120/(ST1\*ST2) MO(2,1)=MO(1,2) С M1(1,1)=C111/(ST1\*ST1) M1(2,2)=C221/(ST2\*ST2) M1(1,2)=C121/(ST1\*ST2) M1(2,1)=C211/(ST1\*ST2) С с С WRITE(6,66) FORMAT(1H1) PRINT OUT THE CORRELATION MATRICES MO AND M1

С WRITE(6,110) ((MO(I,J),J=1,2),I=1,2) WRITE(6,111) ((M1(I,J),J=1,2),I=1,2) FORMAT(5X,'CORRELATION MATRIX: MO',//,((5X,2F10.3)/)) FORMAT(5X,'CORRELATION MATRIX: M1',//,((5X,2F10.3)/)) 110 111 C C C CALCULATE THE PARAMETER MATRICES A AND B CALL LINV2F(M0,2,2,MOINV,0,WKAREA,IER) CALL VMULFF(M1,MOINV,2,2,2,2,2,2,A,2,IER) CALL VMULFP(A,M1,2,2,2,2,2,2,1,2,IER) С DD 10 I=1,2 DD 10 J=1,2 C(I,J)=MO(I,J)-D(I,J) c<sup>10</sup> B(1,1)=C(1,1)\*\*0.5 B(2,1)=C(1,2)/B(1,1) B(2,2)=(C(2,2)-C(1,2)\*\*2/C(1,1))\*\*0.5 B(1,2)=0. CCC PRINT OUT THE MATRICES A AND B WRITE(6,140) ((A(I,J),J=1,2),I=1,2) WRITE(6,141) ((B(I,J),J=1,2),I=1,2) FORMAT(5X,'COEFFICIENT MATRIX: A',//,((5X,2F10.3)/)) FORMAT(5X,'CDEFFICIENT MATRIX: B',//,((5X,2F10.3)/)) 140 141 C 140 NN=N\*12 D0 15 I=1, NN EVR2(I)=VR2(I) EVR1(I)=VR1(I) 15 C C C ESTIMATE THE GAPS OF THE INCOMPLETE SERIES DD 20 I=1,NG I1=ISG(I) I2=IEG(I) K=I2-II+1 DD 40 L=1,K EVR1(I1+L-1)=A(2,1)\*EVR1(I1+L-2)+A(2,2)\*EVR2(I1+L-2) CONTINUE 40 20 0 0 0 0 PERFORM INVERSE TRANSFORMATIONS PP=1/P DD 50 I=1, N DD 50 J=1, 12 L=J+(I-1)\*12 ERAIN1(I,J)=(EVR1(L)\*STD1(J)+XM1(J))\*\*PP ERAIN2(I,J)=(EVR2(L)\*STD2(J)+XM2(J))\*\*PP CONTINUE 50 C C C PRINT OUT THE ESTIMATED SERIES WRITE(6,66) WRITE(6,101) (ID1, (NYEAR(I), (ERAIN1(I,J), J=1,12)), I=1, N) FORMAT(1X,A4,I3,1X,12F6.2) c<sup>101</sup> RETURN 00000000 ----- SUBROUTINE STAT ------SUBROUTINE STAT TRANSFORMS THE ORIGINAL SERIES TO NORMAL AND STATIONARY AND COMPUTES THE STATISTICS OF THE TRANSFORMED SERIES. SUBROUTINE STAT(RAIN, XM, STD, VRAIN, X, ST) DIMENSION RAIN(60, 12), VRAIN(800), XM(12), STD(12) COMMON/A/ ID1, ID2, NYEAR(60) COMMON/B/ N, C,P COMMON/C/ NG, ISG(200), IEG(200) С DO 10 I=1,N DO 10 J=1,12 RAIN(I,J)=(RAIN(I,J)+C)\*\*P CONTINUE 10 CCCC COMPUTE MONTHLY MEANS AND STANDARD DEVIATIONS OF THE NORMALIZED SERIES DO 20 J=1,12 XM(J)=0.

STD(J)=0. DD 25 I=1,N XM(J)=XM(J)+RAIN(I,J)/FLOAT(N) STD(J)=STD(J)+RAIN(I,J)\*\*2 CONTINUE CONTINUE 25 c DD 30 I=1,12 STD(I)=((STD(I)-XM(I)\*\*2\*FLOAT(N))/(FLOAT(N)-1.))\*\*0.5 CONTINUE 30 c c c NOW, STANDARDIZE THE SERIES DO 40 I=1,N DO 40 J=1,12 RAIN(I,J)=(RAIN(I,J)-XM(J))/STD(J) CONTINUE 40 C C C COMPUTE MEAN AND STD OF THE WHOLE SERIES NN=N\*12 IC=0 DD 50 I=1,N DD 50 J=1,12 IC=IC+1 VRAIN(IC)=RAIN(I,J) CONTINUE c<sup>50</sup> X=0. ST=0. D0 60 I=1,NN X=X+VRAIN(I) ST=ST+VRAIN(I)\*\*2 CONTINUE X=X/FLOAT(NN) ST=((ST-X\*\*2\*FLOAT(NN))/(FLOAT(NN)-1.))\*\*0.5 60 X=X/FLOAT(NN) ST=((ST-X\*\*2\*FLOAT(NN))/(FLOAT(NN)-1.))\*\*0.5 C RETURN END //GO.SYSIN DD \* \*\*\*\*\* STATION 6038 - BIVARIATE MODEL \*\*\*\*\* 0 55 25 1 4 27 1 4 27 1 5 3 11 5 33 1 1 5 33 1 1 5 33 2 63 3 1 1 3 63 3 1 1 3 63 3 1 1 3 63 3 1 1 3 63 3 1 1 3 63 3 1 1 3 63 3 1 1 3 7 4 9 2 1 4 2 5 1 2 2 5 1 2 30 2 30 2 //GO.FTI0F001 DD DSN=UF. B0063401. S7. C603820, DISP=(OLD, KEEP) //GO.FT10F001 DD DSN=UF. B0063401. S7. B6093, DISP=(OLD, KEEP) //EDJ С

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