Parallel and Sequential Block Kalman Filtering and Their Implementations Using Systolic Arrays

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Abstract—Two sets of block Kalman filtering equations are derived that differ in the manner of generating the initial and updated estimates. Parallel and sequential schemes for generating these estimates are adopted. It is shown that the parallel implementation inherently leads to a block Kalman estimator which provides filtered estimates at the vector (block) level and fixed-lag smoothed estimates at the sample level. The sequential implementation scheme, on the other hand, generates the estimates of each sample recursively, leading naturally to a scalar (filter) estimator. The scalar estimates are arranged in a vector form, resulting in a block estimator which solely generates filtered estimates both at the vector and sample levels. Simulation results on a speech signal are also presented which indicate the advantages of the sequential block Kalman filter. An algorithm for iterative calculation of Kalman gain and error covariance matrices is given which does not require any matrix inversion operation. The implementation of this algorithm using available systolic array processors is presented. A ring systolic array is also suggested which can be used to implement the state update part of the block Kalman filter.

I. INTRODUCTION

Block processing has been applied to numerous areas in digital signal processing [11]–[19] and control [20]. This is to a great extent due to its advantages and utilities in performing parallel processing with increased throughput rate, increased computational efficiency, reduced roundoff error, and sensitivity performance [11]–[14]. Barnes and Shinnaka [4] used the concept of block processing to arrive at a block state-space formulation with states that propagate only at the edge of each block rather than at each sample. Their block state-space model has been used in a number of papers. In a paper in Anderson et al. [5], this block state-space is used as a dynamic model for modeling the signal and the observation in the Kalman filtering problem. Their block Kalman filter can be used to estimate a block of data at each iteration given the observation blocks up to the present block. Some of the advantages of this scheme, which are inherent to the block processing method are reduced computational effort, and better sensitivity and stability performances. It is also shown that the block estimates are smoother as the block Kalman estimator provides fixed-lag smoothed estimates of the unblocked process. Jain and Jasilek [6] proposed nonrecursive algorithms which utilize FFT to accomplish the operations required in linear smoothing, Riccati equations, boundary value problems, and block Kalman filtering problems. In [7] Lu et al. proposed several systolic architectures for block implemented 1-D FIR and IIR digital filters. These structures offer considerably higher sampling and throughput rates as compared with the single processing element. Parhi and Messerschmitt [8], introduced a look-ahead computation scheme for parallel implementation of the block state-space equation. This scheme utilizes pipelining at bit level and allows an even higher sampling rate. They have also proposed an incremental block state-space structure [9] which offers fewer computational complexities when compared to the standard and parallel block state-space structures.

In this paper two generalized sets of block Kalman filtering equations are derived. It is shown that the manner in which the data is processed determines the types of Kalman filtering equations. In particular, parallel and sequential implementation schemes are considered and the corresponding block Kalman filter equations are obtained. In the first case, the initial estimates are evaluated based upon the updated estimates in the previous block and the estimates are updated in parallel when the entire current data block is received. This scheme is found to be somewhat similar to that developed in [8]. Although the estimates are smoother because of the fixed-lag smoothing property of this structure, the oversmoothing in signal estimation applications may result in loss of some valid information. Moreover, the fixed-lag smoother is sluggish in responding to the fast transitions in the signal. In the second case, the initial estimate for each sample within the block is obtained sequentially from the updated estimates of the previous samples within the same block hence resulting in a more accurate estimation. In addition, this new formulation provides the true filtered estimate of each block without introducing any lag in the estimation. Simulation results on a speech signal are presented which indicate that the sequential block Kalman filter provides the best filtered estimates when compared with those of the original scalar Kalman filter and the parallel block Kalman filter. An iterative procedure for evaluating the Kalman gain and the error covariance matrices is presented which is ideally suited for implementation using current systolic array processors. A ring systolic array is also suggested to implement the state update process involved in any general Kalman filtering algorithm.

II. BLOCK STATE-SPACE FORMULATIONS

Consider a discrete-time, stationary Markov process \( \{ x_k \} \) which is modeled by an autoregressive (AR) model of order \( M \)

\[
x_k = \sum_{i=1}^{M} a_i x_{k-i} + u_k
\]

where \( \{ u_k \} \) represents a white Gaussian (WG) process with zero mean and variance \( \sigma_u^2 \) which drives the autoregression process; and \( a_i \)'s are the coefficients of the AR model. The measurement equation is given by

\[
y_k = \sum_{i=0}^{M} h_i x_{k-i} + v_k
\]
where the sequence \( \{ y_k \} \) represents the observed signal; \( \{ h_k \} \) represents the impulse response sequence of an FIR system of order \( M \); and \( \{ v_k \} \) represents the measurement or observation noise which is assumed to be white with zero mean and variance \( \sigma_v^2 \); and independent of \( \{ u_k \} \), i.e.,

\[
E[u_k v_l] = 0 \quad \forall k, l.
\]

In state-space form the above equations can be written as

\[
S_k = A S_{k-1} + B u_k, \quad E[S_0 S_0^T] = Q_0
\]

\[
y_k = H S_k + v_k
\]

where

\[
S_k = [x_{k-M+1} \cdot x_{k-M+2} \cdots x_k]^T
\]

\[
A = \begin{bmatrix}
0 & 1 & 0 & \cdots & 0 \\
0 & 0 & 1 & \cdots & 0 \\
\vdots & \vdots & \vdots & \ddots & \vdots \\
0 & 0 & 0 & \cdots & 1 \\
a_M & a_{M-1} & \cdots & a_1 & 0
\end{bmatrix}, 
B = \begin{bmatrix}
0 \\
0 \\
\vdots \\
0 \\
1
\end{bmatrix}
\]

\[
H = [h_{M-1} \cdots h_0]
\]

and \( S_0 \) is a zero-mean Gaussian random vector independent of \( \{ u_k \} \) and \( \{ v_k \} \).

Now let us assume that the measurements are received in nonoverlapping blocks of \( N \) samples (\( N = M \) for simplicity), and it is desirable to compute a block of filtered estimates at each stage of the algorithm. The block state-space equation can be obtained by partitioning the sequences \( \{ x_k \}, \{ u_k \}, \{ y_k \}, \) and \( \{ v_k \} \) into nonoverlapping blocks of size \( M \), letting \( k = iM + j \) and writing the state-space equation (4) for all \( j \in [0, M - 1] \). This would yield

\[
X_i = \tilde{A} X_{i-1} + BU_i
\]

\[
Y_i = RX_{i-1} + DU_i + V_i
\]

where

\[
\tilde{A} = A^M
\]

\[
\tilde{B} = \begin{bmatrix}
A^{M-1}B & A^{M-2}B & \cdots & B
\end{bmatrix}
\]

\[
\tilde{H} = \begin{bmatrix}
HA \\
HA^2 \\
\vdots \\
HA^{M-1}
\end{bmatrix}, \quad \tilde{D} = H \tilde{B}
\]

\[
H_e = \begin{bmatrix}
h_0 \\
h_1 & h_0 \\
\vdots \\
h_{M-1} & \cdots & h_0
\end{bmatrix}
\]

and \( X_i \) represents the \( i \)th block of signal \( x_k \), i.e.,

\[
X_i = S_{iM-M+1} \cdot [x_{iM} \cdot x_{iM+1} \cdots x_{iM+M-1}]^T
\]

\( U_i, Y_i, \) and \( V_i \) are similarly defined.

In what follows, two sets of block Kalman filtering equations are derived which differ in the manner of generating the initial \( (a \text{ priori}) \) and the updated \( (a \text{ posteriori}) \) estimates.

### III. PARALLEL BLOCK KALMAN FILTER

In this scheme, the initial estimates of the samples within block “i” are obtained based on the updated estimates within the previous blocks. The updating for the estimates in block “i” will take place when the entire observation block \( Y_i \) is received in parallel. Let us denote the updated \( (a \text{ posteriori}) \) block estimate of \( X_i \), conditioned on the observation blocks up to \( Y_i \) (or, equivalently, observation points up to \( y_{iM+M-1} \)) by \( \hat{X}_i \), i.e.,

\[
\hat{X}_i \triangleq E[X_i | Y_0 \cdots Y_i] = E[X_i | Y_0 \cdots y_{iM+M-1}]
\]

and the initial \( (a \text{ priori}) \) block estimate of \( X_i \), conditioned on the past observation blocks up to \( Y_{i-1} \) (or equivalently observation points up to \( y_{iM-1} \)) by \( \hat{X}_{i-1} \), i.e.,

\[
\hat{X}_{i-1} \triangleq E[X_i | Y_0 \cdots Y_{i-1}] = E[X_i | Y_0 \cdots y_{iM-1}].
\]

Then, the “block innovation sequence” \( \{ Z_i \} \) can be defined as

\[
Z_i = Y_i - H \hat{X}_{i-1}.
\]

Note that in spite of the fact that \( Y_i \) is dependent of \( \{ Y_0 \cdots Y_{i-1} \} \), since \( X_{i-1} \) is obtained based upon the observation set \( \{ Y_0 \cdots Y_{i-1} \} \), the block innovation sequence \( Z_i \) will contain only the “new information” in \( Y_i \) and thus is orthogonal to \( \{ Y_0 \cdots Y_{i-1} \} \). Now, utilizing the linearity property of the minimum variance estimator (MVE) we get

\[
\hat{X}_i = \hat{X}_{i-1} + E[X_i | Z_i]
\]

\[
= \hat{X}_{i-1} + E[X_i | z_{iM} \cdots z_{iM+M-1}]
\]

where the elements of \( Z_i \), i.e., \( z_{iM+j} \) are given by

\[
z_{iM+j} = y_{iM+j} - H \hat{X}_{i-1}
\]

\[
= y_{iM+j} - HA^{i-1} \hat{X}_{i-1}, \quad j \in [0, M - 1].
\]

The initial estimate for each element of a block is obtained recursively based upon observations up to \( y_{iM-1} \), i.e.,

\[
\hat{S}_{iM+j} = [\hat{x}_{(i-1)M+j} \cdots \hat{x}_{iM-1} \hat{x}_{iM} \cdots \hat{x}_{iM+j}]^T
\]

\[
= A \hat{S}_{iM-1} + A^i \hat{X}_{i-1}, \quad j \in [0, M - 2].
\]

For \( j = M - 1 \)

\[
\hat{S}_{iM-1} = [\hat{x}_{iM} \cdots \hat{x}_{iM+M-1}]^T
\]

\[
= \hat{X}_i.
\]

Thus we obtain

\[
\hat{X}_i = A \hat{X}_{i-1} - 1.
\]

Remark III-1: Note that since \( \hat{S}_{iM+j}, j \in [0, M - 1] \) is obtained based upon the observation points up to \( y_{iM-1} \), \( Z_{iM+j} \) does not solely represent the “new part” of \( y_{iM+j} \). Thus

\[
E[z_{iM+k} z_{(iM+1)M+k}^T] \neq 0 \quad k \neq l \quad k, l \in [0, M - 1]
\]

whereas

\[
E[z_{iM+k} z_{(iM+1)M+k-1}^T] = 0 \quad \forall j \in [0, M - 1] \quad \forall l \geq 1.
\]

This fact indicates that although the block innovation sequence \( \{ Z_i \} \) represents a vectorially uncorrelated process, the elements of \( Z_i \) are mutually correlated. In other words, \( Z_i \) can be regarded as a multichannel process [11] with mutual coupling between the individual channels. This contradicts with the scalar nature of the original signal and observation models.
Remark III.2: The block estimate \( \hat{X}_i \) is a vector containing the smoothed estimates of the scalar signal \( x_j \) having different lags depending on the position of the particular element, i.e.,
\[
\hat{X}_i = \begin{bmatrix}
E[x_{iM} | y_0 \cdots y_{iM+M-1}]

E[x_{iM+1} | y_0 \cdots y_{iM+M-1}]

\vdots

E[x_{iM+M-1} | y_0 \cdots y_{iM+M-1}]
\end{bmatrix}
\]
\[
= \begin{bmatrix}
E[x_{iM} | y_0 \cdots y_{iM-1}]

\vdots

E[x_{iM+M-1} | y_0 \cdots y_{iM-1}]
\end{bmatrix}
+ \begin{bmatrix}
E[x_{iM} | Z_i]

\vdots

E[x_{iM+M-1} | Z_i]
\end{bmatrix}
\tag{12}
\]

Using the orthogonal projection lemma [12] and considering that the updating takes place based upon the entire vector \( Z_i \), we have
\[
E[X_i | Z_i] = E[X_i | Z_i] E[Z_i | Z_i]^{-1} Z_i.
\tag{13}
\]

Defining
\[
K(i) \triangleq E[X_i | Z_i] E[Z_i | Z_i]^{-1}
\tag{14}
\]
as the "Kalman gain matrix," the updating equation (8a) becomes
\[
\hat{X}_i = \hat{X}_i + K(i) Z_i.
\tag{15}
\]

New Kalman filter equations are derived which account for the presence of the direct feedthrough term with gain \( D \) in (5). These formulations that differ in complexity from those obtained in [5] are given in order as
\[
\hat{X}_i = \hat{X}_{i-1}
\tag{16a}
\]
\[
R(i) \triangleq E[Z_i | Z_i] = \hat{H}P_a(i-1)\hat{H}^T + \hat{D}Q_D \hat{D}^T + Q_V
\tag{16b}
\]
\[
K(i) = (\hat{H}P_a(i-1)\hat{H}^T + \hat{D}Q_D \hat{D}^T)R(i)^{-1}
\tag{16c}
\]
\[
\hat{X}_i = \hat{X}_i + K(i)[Y_i - \hat{H}\hat{X}_{i-1}]
\tag{16d}
\]
\[
P_a(i) \triangleq E[\hat{X}_{i+1} \hat{X}_i^T] = \hat{H}P_a(i-1)\hat{H}^T + \hat{D}Q_D \hat{D}^T + P_a(i)
\tag{16e}
\]
\[
P_o(i) \triangleq E[\hat{X}_i \hat{X}_i^T] = P_o(i) - K(i)[\hat{H}P_a(i-1)\hat{H}^T + \hat{D}Q_D \hat{D}^T]
\tag{16f}
\]

where \( \hat{X}_i \triangleq X_i \hat{X}_i \), and \( \hat{X}_i \triangleq X_i \hat{X}_i \); \( P_a(i) \) and \( P_o(i) \) are, respectively, the a posteriori and the a priori estimate covariance matrices; \( Q_V \) and \( Q_o \) are the covariance matrices of \{ \hat{U}_j \} and \{ \hat{V}_j \} block sequences, given by
\[
E[U_j U_j^T] = Q_V \delta(i-j) = \sigma_p^2 \delta(i-j)
\]
\[
E[V_j V_j^T] = Q_o \delta(i-j) = \sigma_q^2 \delta(i-j)
\tag{17}
\]

and \( \delta(\cdot) \) represents the Kronecker delta function.

IV. SEQUENTIAL BLOCK KALMAN FILTER

This case differs from the previous one in a number of ways. First, the initial estimate for each sample within a block is formed based upon the updated estimates of all the past samples including the ones in the same block. For example, in finding \( \hat{x}_{iM+j} \), we use all the updated estimates \( \hat{x}_i \), \( iM + j \), \( \cdots \), \( iM + j - 1 \), that are better than \( \hat{x}_i \) used in the previous case. Thus, the estimates generated using this structure are more reliable and accurate. Second, the updating in this case takes place sequentially on each sample within a block as opposed to the parallel method where the updating is done vectorwise. Each processor estimates based upon the relevant observation point available to it, e.g., first processor estimates based upon the first element of the observation block; second processor estimates based upon the first and second observation points, etc. In contrast to the parallel method this scheme generates purely filtered estimates both at the vector and the scalar levels. The updated block estimate in this case is given by
\[
\hat{X}_{iM+j} = \begin{bmatrix}
E[x_{iM} | y_0 \cdots y_{iM-1}]

E[x_{iM+1} | y_0 \cdots y_{iM-1}]

\vdots

E[x_{iM+M-1} | y_0 \cdots y_{iM-1}]
\end{bmatrix}
\tag{18}
\]

which represents the scalar estimates arranged in a block form. The initial block estimate is then given by
\[
\hat{X}_{iM-1} = \begin{bmatrix}
E[x_{iM} | y_0 \cdots y_{iM-1}]

E[x_{iM+1} | y_0 \cdots y_{iM-1}]

\vdots

E[x_{iM+M-1} | y_0 \cdots y_{iM-1}]
\end{bmatrix}
\tag{19}
\]

Applying the linearity property of MVE to each element yields
\[
\hat{X}_i = \hat{X}_i + \begin{bmatrix}
E[x_{iM} | z_{iM}]

E[x_{iM+1} | z_{iM+1}]

\vdots

E[x_{iM+M-1} | z_{iM+M-1}]
\end{bmatrix}
\tag{20a}
\]

where scalar innovation sequence is
\[
\frac{z_{iM+j}}{y_{iM+j} - H \hat{S}_{iM+j}} \quad j \in [0, M-1].
\tag{20b}
\]

Note that in this case \( \hat{S}_{iM+j} \) is given by
\[
\hat{S}_{iM+j} = [\hat{x}_{(i-1)M+j+1} \cdots \hat{x}_{iM+j}] R_{iM+j}^{-1} [\hat{x}_{iM+j} \cdots \hat{x}_{(i-1)M+j+1}]
\tag{21}
\]

Also we have
\[
\hat{S}_{iM-1} = \hat{X}_{i-1}
\tag{22a}
\]

and
\[
\hat{S}_{iM+M-1} = \hat{X}_i
\tag{22b}
\]

Remark IV.1: In this case \( \hat{S}_{iM+j}, j \in [0, M-1], \) in the scalar innovation is obtained based upon the updated estimates \( \hat{x}_i, i = iM + j, \cdots, iM + j - 1 \), which are in turn evaluated based upon the observation points up to \( y_{iM+j-1} \). That is, \( z_{iM+j} \) does indeed represent the "new part" of \( y_{iM+j} \), and thus is independent of all the observation points in the set \( \{ y_0 \cdots y_{iM+j-1} \} \), i.e.,
\[
E[z_{iM+j} | y_i] = 0 \quad \forall i \in [0, iM + j - 1]
and
\[ E[z_k z_l] = 0 \quad k \neq l, \quad \forall k, l \]  
(23)

This implies that the sequence \{z_k\} represents an uncorrelated scalar process. This certainly matches more closely to the scalar nature of the original signal and observation models.

Applying the linearity property of the estimator to each scalar estimate yields
\[ E[S_{m+j} | y_0 \cdots y_{m+j}] = E[S_{m+j} | y_0 \cdots y_{m-1}] + E[S_{m+j} | z_{m+j}] \]

or
\[ \hat{s}_{m+j} = S_{m+j} + E[S_{m+j} | z_{m+j}] \]  
(24)

Now using the orthogonal projection theorem [12] the second term on the right-hand side of (24) can be expressed as
\[ E[S_{m+j} | z_{m+j}] = E[S_{m+j} | z_{m+j}] \left( E[z_{m+j} z_{m+j}]^{-1} \right)^{-1} \]

\[ = K(j) z_{m+j} \]  
(25)

where \( K(j) \) is the Kalman gain matrix for the jth element in block "i". The Kalman filter equations for this structure are derived to be
\[ \hat{s}_{m+j} = \hat{s}_{m+j-1} \quad j \in [0, M - 1] \]  
(26a)

\[ P_a(iM + j) = A P_a(iM + j - 1) A^T + B B^T \sigma_e^2 \]  
(26b)

\[ K(j) = P_a(iM + j) H^T (H P_a(iM + j) H^T + \sigma_e^2)^{-1} \]  
(26c)

\[ z_{m+j} = y_{m+j} - H \hat{s}_{m+j}, \quad j \in [0, M - 1] \]  
(26d)

\[ \hat{s}_{m+j} = \hat{s}_{m+j} + K(j) z_{m+j} \]  
(26e)

\[ P_a(iM + j) = [I - K(j) H] P_a(iM + j) \]  
(26f)

Now writing (26e) for \( j \in [0, M - 1] \), i.e., for each element of block "i" and expressing all the initial estimate vectors in terms of \( \hat{x}_{-1} \) yields
\[ \hat{s}_{m-M-1} = \hat{x}_{-1} = A^M \hat{x}_{-1} + K z_{i} \]  
(27)

where
\[ K_i = \left[ A^{M-1} K_0 \cdots K(M-1) \right]. \]

Similarly writing (26d) in block form and using (21) and (22) gives
\[ Z_i = Y_i - H \hat{x}_{-1} - \hat{R}_i Z_i \]  
(28)

where \( \hat{H} \) is defined in (5) and \( \hat{R}_i \) is given by its \((k, l)\) elements as
\[ \hat{R}_i(k, l) = \begin{cases} 0 & k \leq l \\ H a^{k-l} & k > l \end{cases} \quad k, l \in [1, M]. \]

Equation (28) can be rewritten as
\[ Z_i = \hat{W}_i^{-1} (Y_i - \hat{H} \hat{x}_{-1}) \]  
(29a)

where the nonsingular matrix \( \hat{W}_i \) is
\[ \hat{W}_i = I + \hat{R}_i. \]  
(29b)

Equations (27) and (29) together with (26b), (26c), and (26f) for calculating Kalman gain and error covariances represent a set of sequential block Kalman filter equations, obtained by blocking the scalar estimator equations. As can be seen, these formulations provide purely block filtered estimate. If stationarity is assumed, the steady-state Kalman gain \( \hat{K} \) can be obtained using (26b), (26c), and (26f). This gain matrix is then used throughout the entire updating process. Equations (27) and (29) would then represent a simplified sequential block Kalman filtering procedure.

Remark IV.2: In addition to the above benefits, this scheme provides a very efficient means of block Kalman filtering, since the Kalman gain and error covariance evaluations using (26b), (26c), and (26f) do not require any matrix inversion operation, as in the scalar case.

V. SIMULATION RESULTS

In order to examine the effectiveness of the proposed block Kalman filters, a speech signal for the word "hello" is digitized at 16 kHz and used throughout the experiment. This signal which contains 5000 points is shown in Fig. 1. This signal is then corrupted by adding white Gaussian noise with zero mean and variance 0.158 to obtain a SNR of 6.9 dB. The corrupted signal is shown in Fig. 2. The entire simulation is performed on a Mac II computer with a 68020 processor and a 68881 Math coprocessor. A fifth-order AR model is fitted to this data and the model parameters are obtained by solving the Yule–Waker equations. Using this method the variance of the driving noise sequence \( \sigma_e^2 \) is found to be 0.02. Standard Kalman filter is first applied to the scalar model in (4) with initial \( P(0) = 10 I \) and \( H = 10 0 0 \cdots 0 1 \). The filtered signal is shown in Fig. 3. The SNR is measured to be 9.7 dB which shows considerable noise reduction. The entire process is implemented on line and it took 148 s. Fig. 4 shows the behavior of the parallel block Kalman filter. The effect of smoothing can clearly be seen in this result. This filter has slower response than the standard scalar Kalman filter. The degradation in performance can also be observed from the variance of the output noise that is found to be 0.286, giving a SNR of 4.3 dB which is even smaller than that of the degraded signal. Additionally, some less of valid information is evident in the processed signal. The computational time for this process is approximately 94 s which is smaller than that of the standard Kalman filter. Again in this case the Kalman gain equations are evaluated on-line. The result of the sequential block Kalman filter is shown in Fig. 5 which indicates the effectiveness of this method when compared with the previous cases. This improvement is also reflected in the value of the variance of the noise in the output signal which is reduced to 0.038 giving a SNR of 11.25 dB. The estimates are obtained in approximately 31 s. In this case the steady state Kalman gain was obtained off-line, thus (27) and (29) are the only ones that are evaluated on-line. Overall, the tracking behavior of the sequential block Kalman filter is by far better than the parallel case. This, coupled with its accuracy and efficiency as discussed above makes this method very attractive for signal estimation applications.

VI. IMPLEMENTATION USING SYSTOLIC ARRAYS

The development of Kalman filtering algorithms suitable for implementation using systolic arrays has recently been considered in a number of papers, namely [13]–[15]. In this section, systolic array implementations of the block Kalman filter are presented. The results can be applied to any general MIMO Kalman filter structure. The proposed systolic architectures im-
implement all the operations involved in the Kalman filtering process such as the Kalman gain evaluation, iterative update for the a priori and the a posteriori error covariance matrices and the state update procedure. An algorithm for iterative calculations of Kalman gain and the error covariance matrices is suggested which does not require any matrix inversion operation. The global convergence of this algorithm is also established. The consecutive matrix-matrix multiplications (MMM’s) in this algorithm are implemented on an expanded skew pipeline structure which uses the available geometric array parallel processor (GAPP) [15], [16] type systolic arrays. The state update process is implemented using a ring systolic array structure [17]-[19]. Let us first briefly introduce the GAPP systolic array.

**A. GAPP [16]**

The geometric array parallel processor or GAPP is a commercial two-dimensional systolic array processor chip manufac-
tured by NCR. It is a mesh-connected 6-by-12 arrangement of 1-b processor elements, or PE's. Each PE can communicate with four neighbors: North, East, South, and West. Each PE is composed of a 1-b serial ALU, 128 word by 1-b RAM and 4 single bit latches: three latches hold inputs to the ALU and the fourth latch allows I/O through the cell without interrupting the ALU, i.e., I/O operations are overlapped with the computations. Each instruction is broadcast to all the PE's making the array to operate like a single instruction multiple data (SIMD) machine.

The cascadeability of the GAPP allows system designers to implement arrays of processors of arbitrary size in multiples of 6-by-12 elements. Fig. 6 shows the interconnections among four cells. The following development is based upon a generalized GAPP type systolic array and not limited by the specific size of memory and/or the dimension.

**B. Consecutive Matrix-Matrix Multiplication Using GAPP**

A GAPP type array is shown in Fig. 7(a). External connections are added to this array in order to perform the operations needed in this section. The resultant array is shown in Fig. 7(b). Note that the required multiplexers for the external communication links are not shown in this figure for the sake of simplicity.

Let us assume that a matrix, say $A$, is cyclically loaded from the bottom of the array in a row parallel fashion as shown in Fig. 8(a). The matrix stored in the array shall be denoted by $A^*$. Note that the diagonal shift operations in this array can be accomplished in two steps. Fig. 8(b) shows the loading of a $3 \times 3$ matrix. Once the loading operation is completed, all the MMM combinations of matrix $A$ (or $A'$) by matrix $B$ (or $B'$) which result in $C$ (or $C'$) can be computed in a skew pipelined fashion as shown in Fig. 9. Without loss of generality and to simplify the discussion here we have assumed that matrices $A$, $B$, and $C$ are of the same size as $M \times M$. At every PE, an entry of matrix $B$ is multiplied by the corresponding entry of matrix $A$ and the result is added to the entry of matrix $C$. Then all the entries of matrices $B$ and $C$ move to other PE's following their corresponding paths. All the possible combinations of MMM's and their data flows are shown in Table I. A $3 \times 3$ example for performing matrix operation $BA$ and the types of inputting and outputting are shown in Fig. 10. As can be seen, the partial summations for matrix $C$ is skewed for this particular operation.

For a GAPP type SIMD machine, one cannot change the stored matrix $A^*$ before the completion of the first MMM (before the last row or column of $C$ is generated). Thus, the efficiency of this structure is limited only to 50% if the array is to be used for consecutive MMM operations. However, $C$ or $C^*$ can be loaded into the same array, but with different memory
address, during the MMM time for the next consecutive MMM operation as shown in Fig. 11. If the type of inputting of the next MMM matches the type of outputting of the present MMM (see Table I), e.g., both are row parallel, then the two MMM’s can be performed using a pair of identical arrays without interruption. The relevant architecture which improves the speed by a factor of two is shown in Fig. 12. These architectures will be used to implement the consecutive MMM operations involved in the iterative update of Kalman gain and also the error covariance matrices.

In what follows, an algorithm for iterative computation of Kalman gain and error covariance matrices is proposed which avoids any matrix inversion operation.

### Table I

<table>
<thead>
<tr>
<th>Type of Inputting B</th>
<th>Path of B</th>
<th>Summing Path of C</th>
<th>Type of Outputting C</th>
</tr>
</thead>
<tbody>
<tr>
<td>A B</td>
<td>Parallel</td>
<td>Skewed</td>
<td>Column</td>
</tr>
<tr>
<td>A' B</td>
<td>Parallel</td>
<td>Left-to-right</td>
<td>Skewed</td>
</tr>
<tr>
<td>B A</td>
<td>Parallel</td>
<td>Left-to-right</td>
<td>Row</td>
</tr>
<tr>
<td>B A</td>
<td>Parallel</td>
<td>Left-to-right</td>
<td>Row</td>
</tr>
<tr>
<td>B A'</td>
<td>Parallel</td>
<td>Skewed</td>
<td>Row</td>
</tr>
<tr>
<td>B A'</td>
<td>Parallel</td>
<td>Skewed</td>
<td>Row</td>
</tr>
<tr>
<td>A B</td>
<td>Parallel</td>
<td>Skewed</td>
<td>Column</td>
</tr>
<tr>
<td>A B</td>
<td>Parallel</td>
<td>Skewed</td>
<td>Column</td>
</tr>
</tbody>
</table>

### Figure 11

Recall from (16b) that

\[
R_i(i) = \bar{H}P_a(i - 1)\bar{H}^T + \bar{D}Q_0\bar{D}^T + Q_i.
\]

In this expression all the terms except \(P_a(i - 1)\) are constants. If the Kalman filter is assumed to be convergent (the convergence issue will be discussed later), then the elements of \(P_a\) matrix decrease from one iteration to the next. Thus we can write

\[
R_i(i) = R_i(i - 1) - \Delta R_i(i - 1)
\]

where the error matrix \(\Delta R_i(i - 1)\) is

\[
\Delta R_i(i - 1) = \bar{H}P_a(i - 2) - P_a(i - 1)\bar{H}^T
\]

\[
= \bar{H}\Delta P_a(i - 1)\bar{H}^T.
\]

The inverse of \(R_i(i\) can also be written as

\[
R_i^{-1}(i) = R_i^{-1}(i - 1) + \Delta R_i^{-1}(i - 1).
\]

Note that \(\Delta R_i^{-1}(i - 1) = \Delta R_i^{-1}(i - 1). Premultiplying (32) by (31a) gives \n\Delta R_i^{-1}(i - 1) = R_i^{-1}(i - 1) - \Delta R_i(i - 1)R_i^{-1}(i - 1)
\[
+ \Delta R_i^{-1}(i - 1)\Delta R_i(i - 1)R_i^{-1}(i - 1).
\]

If we ignore the effects of the second term on the right-hand side of (33) the algorithm in [20] will be obtained. Now let us denote \(R_i^{-1}\) by \(\Phi\) and \(\Delta R_i^{-1}\) by \(\Delta \Phi\), then the following iterative
procedure for solving (33) can be used:

$$\Delta \Phi_i(i - 1) = \Phi(i - 1) \Delta R_i(i - 1) \Phi(i - 1)$$

where \( \Delta \Phi_i(i - 1) \) is the \((k + 1)\)th order solution of (33) at the \(i\)th iteration of the filtering algorithm. For \(k = 0\), the first-order approximation is:

$$\Delta \Phi_0(i - 1) = \Phi(i - 1) \Delta R_0(i - 1) \Phi(i - 1)$$

which is the solution used in [20].

Expression (34) can be rewritten as:

$$\Delta \Phi_i(i - 1) = \Phi_{k - 1}(i) \Delta R_k(i - 1) \Phi(i - 1) \quad \forall k \geq 0$$

(36a)

where:

$$\Phi_{k - 1}(i) = \Phi(i - 1) + \Delta \Phi_{k - 1}(i - 1)$$

(36b)

Note that \( \Phi_{k - 1}(i) \) is the \(k\)th order estimate of \( \Phi(i) \). Using (36) we have:

$$\Phi_k(i) = \Phi(i - 1) + \Delta \Phi_k(i - 1)$$

$$= [I + \Phi_{k - 1}(i) \Delta R_k(i - 1)] \Phi(i - 1), \quad \forall k \geq 0.$$  

(37)

This offers a recursive equation for generating the estimate of \(R_k(i)\) (or \(\Phi(i)\)) with arbitrary accuracy given the matrices \(\Delta R_0(i - 1)\) and \(\Phi(i - 1)\). The error matrix \(\Delta R_k(i - 1)\) is computed from (31b) and \(\Phi(i - 1)\) is the final solution of \(\Phi\) (or \(R_k^{-1}\)) at filtering iteration \((i - 1)\).

**Lemma VI.1:** The recursive algorithm in (37) is globally convergent in the sense that

$$\lim_{k \to \infty} \| \Phi_k(i) - \Phi_{k - 1}(i) \| = 0.$$  

(38)

**Proof:** From (37) we can write:

$$\Phi_k(i) - \Phi_{k - 1}(i)$$

$$= [\Phi_{k - 1}(i) - \Phi_{k - 2}(i)] \Delta R_{k - 1}(i - 1) \Phi(i - 1)$$

$$= [\Phi_{k - 1}(i) - \Phi_{k - 2}(i)] \Delta R_{k - 1}(i - 1) \Phi(i - 1)$$

(39)

Now, using (34)-(36) yields:

$$\Phi_k(i) - \Phi_{k - 1}(i)$$

$$= \Phi(i - 1) \Delta R_k(i - 1) \Phi(i - 1)^{k - 1}.$$  

(40)

Taking the norm of both sides of (40) gives the following inequality:

$$\| \Phi_k(i) - \Phi_{k - 1}(i) \| \leq \| \Phi(i - 1) \| \| \Delta R_k(i - 1) \| \Phi(i - 1) \|^{k - 1}$$

$$\leq \lambda_{\text{min}}[R(i - 1)] \| \Delta R_k(i - 1) \| \Phi(i - 1) \|^{k - 1}$$  

(41)

where \(\lambda_{\text{min}}[R(i - 1)]\) represents the smallest eigenvalue of \(R_k(i - 1)\). The second term on the right-hand side is \(\Delta R_k(i - 1)\Phi(i - 1)\):

$$= (\Phi \Delta P_a(i - 1) \Phi^T) [ \Phi \Delta P_a(i - 1) \Phi^T + \Delta Q_a \Phi \Phi^T + Q_a ]^{-1}.$$  

(42)

Thus it is clear that

$$\| \Delta R_k(i - 1) \Phi(i - 1) \| \leq 1.$$  

Additionally, due to the presence of \(Q_r \) and \(\Delta Q_r \Phi \Phi^T \) terms in \(R_k(i - 1)\):

$$\lambda_{\text{min}}[R(i - 1)] \leq 1.$$  

Thus

$$\lim_{k \to \infty} \| \Phi_k(i) - \Phi_{k - 1}(i) \| = 0.$$  

Q.E.D.

As a result, the iterative procedure for on-line computation of the Kalman gain and error covariance matrices using the above algorithm becomes:

**Step 1:** Compute the error matrix

$$\Delta R_k(i - 1) = \Phi \Delta P_a(i - 2) \Phi^T + \Delta Q_r \Phi \Phi^T + Q_r.$$  

(43a)

**Step 2:** Iterate

$$\Phi_k(i) = [I + \Phi_{k - 1}(i) \Delta R_k(i - 1)] \Phi(i - 1), \quad \forall k \geq 0$$  

(43b)

with

$$\Phi_{k - 1}(i) = \Phi(i - 1)$$  

(43c)

for \(k = 0, 1, \ldots, \) until for \(k = k^*\) (say) we have

$$\Phi_k(i) \Delta R_k(i - 1) \Phi(i - 1) \prec \epsilon.$$  

\(\forall m, n[1, M]$$

where \(\Phi_k(i) \Delta R_k(i - 1) \Phi(i - 1)\) represents \((m, n)\)th entry of \(\Phi_k(i)\) matrix and \(\epsilon\) is a small positive threshold. Then set \(\Phi(i) = \Phi_k(i)\) and proceed.

**Step 3:** Evaluate Kalman gain and error covariances using

$$K(i) = (\Phi \Delta P_a(i - 1) \Phi^T + \Delta Q_r \Phi \Phi^T) \Phi(i - 1)$$

(43d)

$$P_a(i) = \Phi \Delta P_a(i - 2) \Phi^T + \Delta Q_r \Phi \Phi^T + Q_r$$

$$- K(i) [ \Phi \Delta P_a(i - 1) \Phi^T + \Delta Q_r \Phi \Phi^T + Q_r ]^{-1}$$

(43e)

Repeat these steps for all \(i\).

This algorithm is very well suited for implementations using the systolic array architectures designed in Section VI-B. This is considered in the next section.

**D. Systolic Array Implementation of Kalman Filter Equations**

Using the structure in Fig. 12 an implementation scheme can be devised which provides minimum computation time when the inputting and outputting patterns are matched in accordance with Table II. The critical (longest) path of successive matrix operations in algorithm (43) when implemented on this structure can be as small as 5.5 MM on this particular case occurs when the first-order approximation of \(\Phi(i)\) would suffice. A M/M time is defined as the time elapsed between the first vector of the input matrix entering the array and the last vector of the output matrix leaving the array. The decision rule in step 2 of the algorithm can be implemented using the global output of the GAPP type architecture [16]. Assuming that \(k^*\) is the value of \(k\) at which the appropriate solution for \(\Phi(i)\) is obtained, then from Table II one iteration period is

$$T = (4k^* + 11)T.$$  

(44a)
TABLE II
SCHEDULE FOR COMPUTING KALMAN GAIN AND ERROR COVARIANCE MATRICES USING THE ARCHITECTURE IN FIG. 12

<table>
<thead>
<tr>
<th>MMM Period</th>
<th>Input Type</th>
<th>Operation</th>
<th>Output Type</th>
<th>Operation</th>
<th>Output Type</th>
</tr>
</thead>
<tbody>
<tr>
<td>T₁</td>
<td>C</td>
<td>L₁ = R*₆P₆(i - 1)</td>
<td>C</td>
<td>M₆ = Ł₆P₆(i - 1)</td>
<td>C</td>
</tr>
<tr>
<td>T₂</td>
<td>R</td>
<td>L₃(i) = L₃(i)R₆'</td>
<td>R</td>
<td>M₆ = M₆R₆'</td>
<td>R</td>
</tr>
<tr>
<td>T₃</td>
<td>R</td>
<td>L₅(i) = L₅(i) - L₅(i - 1)</td>
<td>R</td>
<td>M₆ = M₆ + (BQ₆)D'</td>
<td>R</td>
</tr>
<tr>
<td>T₄</td>
<td>R</td>
<td>L₅(i) - Φ(i - 1)ΔR₅(i)</td>
<td>R</td>
<td>M₆ = M₆Ł₆</td>
<td>R</td>
</tr>
<tr>
<td>T₅</td>
<td>C</td>
<td>L₆ = I + L₆</td>
<td>C</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
<tr>
<td>T₆</td>
<td>C</td>
<td>L₇(i) = Φ(i - 1)ΔR₇(i)</td>
<td>C</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
<tr>
<td>T₇</td>
<td>R</td>
<td>L₉(i) = Φ(i)ΔR₉(i)</td>
<td>R</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
<tr>
<td>T₈</td>
<td>C</td>
<td>L₈(i) = Φ(i)ΔR₈(i)</td>
<td>C</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
<tr>
<td>T₉</td>
<td>R</td>
<td>L₁₀(i) = L₁₀(i) + L₁₀</td>
<td>R</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
<tr>
<td>T₁₀</td>
<td>C</td>
<td>L₁₁(i) = Φ(i - 1)ΔR₁₁(i)</td>
<td>C</td>
<td>M₆ = M₆L₆</td>
<td>C</td>
</tr>
</tbody>
</table>

Where, R = Row parallel, C = Column parallel.

where \( T_i \) is one half of a MMM computation time, i.e.,
\[
T_i = M(T_{MUL} + T_{ADD})
\]
and \( T_{MUL} \) is the scalar multiplication time and \( T_{ADD} \) is the scalar addition time. Note that it is assumed that the results of \( BQ₆D' \) and \( BQ₆D' \) are precomputed and stored in the array. The calculation of the state vector \( \hat{X}_i \) can be inserted during any idle period of the array processor in Fig. 12.

If the Kalman gain matrix is preevaluated off-line on a host computer, the state update part can be implemented using a simpler structure described in the next section.

E. Systolic Array Implementation of State Update Part

In our earlier work [17] several parallel and pipeline algorithms and structures for 2-D recursive filters were proposed. The skew pipeline structure presented in [17] uses a GAPP type array for 2-D scalar and block filtering operations. In this section, a skew pipeline architecture is presented for both parallel and sequential block Kalman filters which uses a 1-D systolic array in which each PE is assigned to the computation of an element within the block estimate. The 1-D skew pipeline structure is quite similar to the ring systolic array in [18] and [19].

In the development of this architecture, it is assumed that the steady-state Kalman gain matrices are evaluated off-line and used throughout the updating process. In this case (16a) and (16d) can be rewritten as
\[
\hat{X}_i = \hat{X}_{i-1} + K_i[Y_i - \hat{\theta}\hat{X}_{i-1}]
\]
(45a)

or
\[
\hat{X}_i = D\hat{X}_{i-1} + K_iY_i
\]
(45b)

where
\[
D \triangleq \hat{A} - K_i\hat{B}
\]
(46)

and \( K_i \) is the steady-state Kalman gain matrix for the parallel block Kalman filter. Similarly, (27) and (29) can be rewritten as
\[
\hat{X}_i = \hat{X}_{i-1} + K_i\hat{W}_i^{-1}(Y_i - \hat{\theta}\hat{X}_{i-1})
\]
(47a)

or
\[
\hat{X}_i = F\hat{X}_{i-1} + GY_i
\]
(47b)

where
\[
D \triangleq \hat{A} - K_i\hat{W}_i^{-1}\hat{B}
\]
\[
G \triangleq \hat{X}_i
\]

and \( \hat{X}_i \) is the steady-state Kalman gain matrix for the sequential block Kalman filter. The processes in (45b) and (47b) can now be implemented using only two matrix-vector multiplications. A single PE is assigned to each element of the block estimate \( \hat{X}_i \). Fig. 13(a) shows the configuration of the 1-D ring systolic array processor, for \( M = 4 \), where \( M \) is the block or the array size. The structures of all PE's are identical. Fig. 13(b) shows the internal structure of \( PE_m \), i.e., the \( m \)th PE where \( \alpha_{m,n} \) and \( \beta_{m,n} \), \( m, n \in [0, M - 1] \), denote the \( m, n \)th elements of the matrices \( D \) (or \( F \)) and \( K_i \) (or \( G \)) in (45b) (or (47b)), respectively. The elements of these matrices in \( PE_m \) are stored in the following order: \{ (m, m), (m, (m + M - 1) mod M), (m, (m + M - 2) mod M), \ldots, (m, (m + 1) mod M) \} \text{ m} \in [0, M - 1].
During each iteration, these coefficients are shifted in a round-robin fashion. $R_m$ and $S_m$ are two registers used to store the variables. It is assumed that before stage $i$, $\hat{x}_{i-1}(m)$, i.e., the $m$th element of vector $\hat{x}_{i-1}$ has been generated and held in register $S_m$. The operations at each iteration start by shifting $\hat{x}_{i-1}(m)$ from register $S_m$ to $R_m$ and clearing $S_m$. Then in the next $2M$ steps the first and second terms in (45b) or (47b) are generated consecutively in registers $S_m$. At the completion of the last step, the elements of the updated estimate vector $\hat{x}_i$ can be obtained from registers $S_m$ to $S_{m+1}$. The process continues until all the block estimates are computed. The average sampling period for each element of a block is

$$T_s = 2(T_{\text{MUL}} + T_{\text{ADD}})$$

where $T_{\text{MUL}}$ and $T_{\text{ADD}}$ are defined before. The delay time (latency) from $y_i$ arrival to $\hat{x}_i$ generation is

$$T_D = M(T_{\text{MUL}} + T_{\text{ADD}}).$$

VII. CONCLUSION

Two different sets of block Kalman filter equations are derived which correspond to the parallel and sequential modes of implementation. It is shown that the parallel implementation provides fixed-lag smoothed estimates. However, in this scheme the estimation cannot start until the entire observation block is available. The sequential block implementation does not have this particular problem. The initial estimates are more accurate in this latter scheme since they are produced based upon the updated estimates of the samples in the previous block and also those in the current block. This scheme provides purely filtered estimates both at the block and scalar levels and also does not introduce any delay (lag) in the response. A comprehensive study among these two structures and the standard Kalman filter is made which reveals the effectiveness of the sequential block Kalman filter for signal estimation applications. The implementation of the block Kalman filter algorithm using systolic array is also considered. An iterative procedure for efficient computation of Kalman gain and error covariance matrices is proposed which is free of any matrix inversion operation. A general systolic array architecture which utilizes a GAPP-type processor is designed to implement these operations. The state update part when the steady-state Kalman gain is evaluated off-line can be implemented using a simple ring systolic array structure.

REFERENCES


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