

7.2 Linear Interference Cancellation

7.2.1

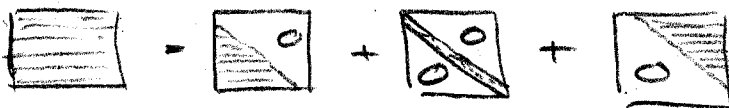
- Linear interference cancellation refers to iterative solution of the ZF or MMSE equations, in which previous estimates of interferers are subtracted (cancelled) out.
- The converged solution is just a ZF or MMSE receiver. We already know their performance (Section 4.4), so why study them further? Because the form of the equations is instructive, and suggests variations can be implemented more easily. [Tan00, R00y00]
- After MF/Rake/MRC reduction to sufficient stats, we want to solve

$$G \underline{b} = \underline{y} \quad \text{where } G = C^T R C A \quad \text{ZF}$$

$$G = (C^T R C + N_0 A^{-2}) A \quad \text{MMSE}$$

Additive decomposition of G helps:

$$G = L + D + U$$



$$\text{want } (L + D + U) \underline{b} = \underline{y} \quad \underline{b} \in \mathbb{R}^{NK}$$

How?

• Jacobi iteration (Appendix 0) is one way:

$$D \underline{b} = \underline{\zeta} - (L+U) \underline{b} \quad \text{is satisfied by a solution.}$$

The j^{th} iterate is $\underline{b}^{(j)}$, and update as

$$D \underline{b}^{(j)} = \underline{\zeta} - (L+U) \underline{b}^{(j-1)}$$

$$\underline{b}^{(j)} = D^{-1} (\underline{\zeta} - (L+U) \underline{b}^{(j-1)})$$

The estimate of bit i (of KU) is determined by previous iteration of all other bits:

$$b_i^{(j)} = D_{ii}^{-1} \left(\zeta_i - \sum_{r \neq i} (L+U)_{i,r} b_r^{(j-1)} \right) \quad i = 1, 2, \dots, NK$$

Convergence iff all eigenvals of $D^{-1}(L+U)$ are in the unit circle.

• Convergence is usually faster with Gauss-Seidel, in which we recognize that the update is typically performed row by row — so use the updated estimates from previous rows of the same iteration

$$b_i^{(j)} = D_{ii}^{-1} \left(\zeta_i - \sum_{r < i} (L+U)_{i,r} b_r^{(j)} - \sum_{r > i} (L+U)_{i,r} b_r^{(j-1)} \right)$$

or

$$\underline{b}^{(j)} = D^{-1} (\underline{\zeta} - L \underline{b}^{(j)} - U \underline{b}^{(j-1)})$$

Convergence iff all eigenvals of $(I + D^{-1}L)^{-1} D^{-1}U$ are in the unit circle, since

$$(I + \bar{D}'L) \underline{b}^{(i)} = -\bar{D}'U \underline{b}^{(i-1)} + \bar{D}'\underline{s}$$

$$\underline{b}^{(i)} = -(I + \bar{D}'L)^{-1} \bar{D}'U \underline{b}^{(i-1)} + (I + \bar{D}'L)^{-1} \bar{D}'\underline{s}$$

- There are about $(NK)^2$ operations in each iteration sweep, so want the number of iterations to be small. We only have to bring the residual error down to a little less than the noise level, so shouldn't take many.

Block Based Formulation

- Neither of the formulations above recognized or exploited the block structure of the equations.

$$G \underline{b} = \underline{s}$$

$$\begin{bmatrix} G_{11} & G_{12} & 0 & 0 & 0 & \dots & 0 \\ G_{21} & G_{22} & G_{23} & 0 & 0 & \dots & 0 \\ 0 & G_{32} & G_{33} & G_{34} & 0 & \dots & 0 \\ \vdots & \vdots & \vdots & \vdots & \vdots & \ddots & \vdots \\ 0 & 0 & 0 & \dots & G_{N-1,N} & \dots & 0 \\ 0 & 0 & 0 & \dots & G_{N,N} & \dots & 0 \end{bmatrix} \begin{bmatrix} \underline{b}(1) \\ \underline{b}(2) \\ \vdots \\ \underline{b}(N) \end{bmatrix} = \begin{bmatrix} \underline{s}(1) \\ \underline{s}(2) \\ \vdots \\ \underline{s}(N) \end{bmatrix}$$

It's organized by $K \times K$ blocks. Organize the iteration in the same way.

- Define the block-based additive decomposition of G :

$$G = L_B + D_B + U_B$$

We won't use the tridiagonal structure just yet. For the moment, just express the iteration in block form as follows:

Jacobi

$$D_B \underline{b}^{(j)} = \underline{\zeta} - (L_B + U_B) \underline{b}^{(j-1)} \quad \text{as before}$$

For symbol time n ,

$$D_{B,n,n} \underline{b}^{(j)} = \underline{\zeta}^{(n)} - \sum_{r \neq n} (L_B + U_B)_{n,r} \underline{b}^{(j-1)}$$

Gauss Seidel

$$D_{B,n,n} \underline{b}^{(j)} = \underline{\zeta}^{(n)} - \sum_{r < n} L_{B,n,r} \underline{b}^{(j)} - \sum_{r > n} U_{B,n,r} \underline{b}^{(j-1)}$$

Now we have a matrix to invert ($D_{B,n,n}^{-1}$) for each time step, so it seems not much point to this approach. However, it can also be solved iteratively (an "inner iteration").

• We now have hybrid approaches

outer iteration

inner iteration

Jacobi

Jacobi

Gauss Seidel

Gauss-Seidel

other (e.g. conjugate gradient)

The inner iteration can be carried out to effective convergence or left after a few iterations.

• The computation is determined by the number of outer and inner iterations J_{out}, J_{in} . Total computation

$$J_{out} N (J_{in} + N) K^2 \rightarrow J_{out} N (J_{in} + 2) K^2$$

recognizing tridiagonal

• How does performance depend on number of iterations?

Explicit matrix inversion for inner

Allows block size to differ from K

Asynch. Flat AWGN

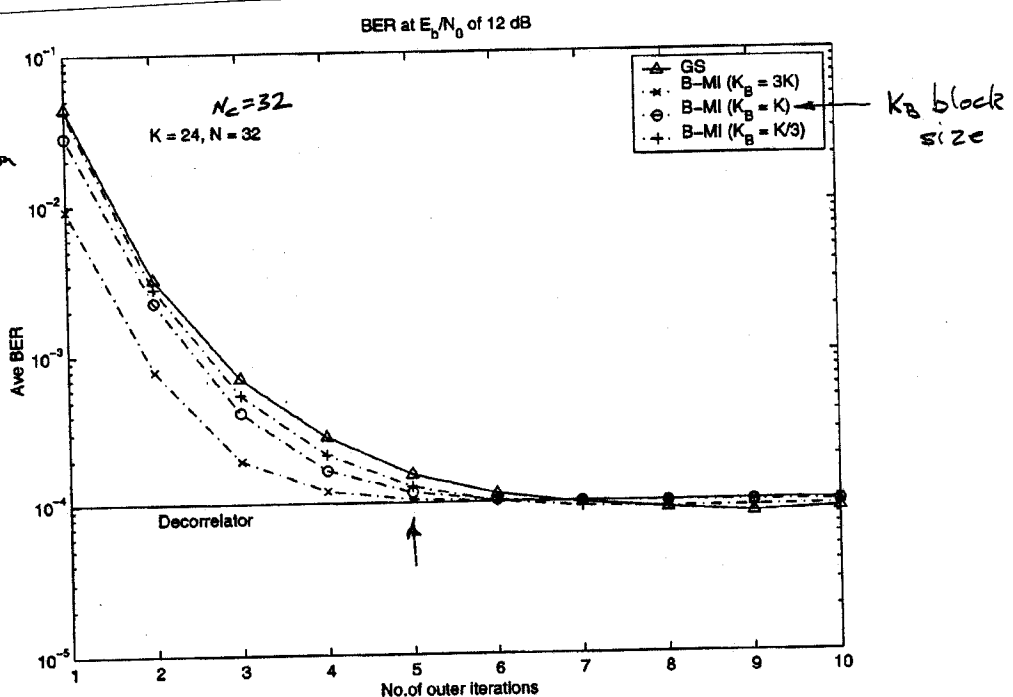
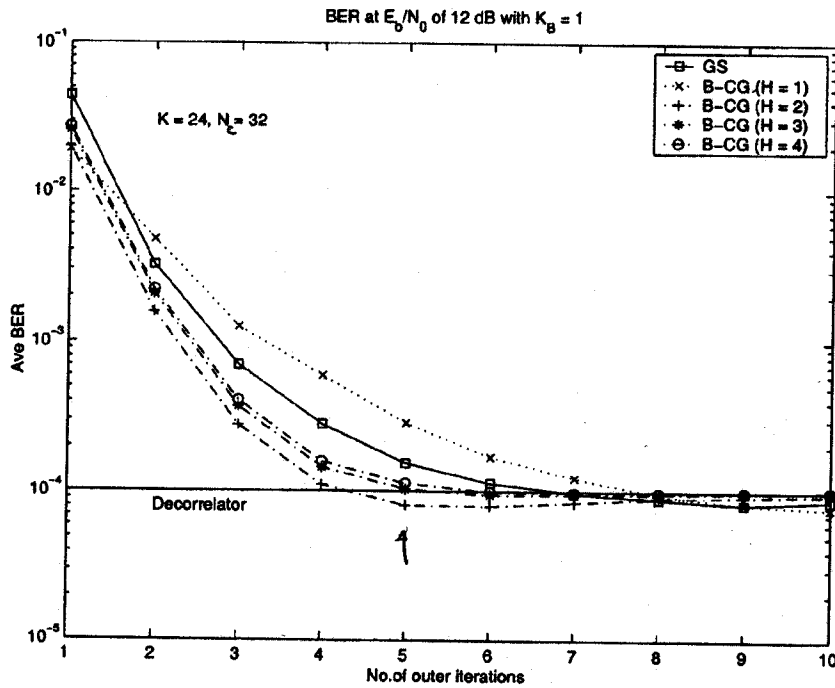


Fig. 5. Averaged BER of the B-MI. Block sizes of $K_B = 8, 24, 72$ are considered where $N = 32$ and $E_b/N_0 = 12$ dB. The point GS is included for reference.

$N_c = 32$

Conjugate
gradient inner
iteration.



$H = J_{in}$

CG is best (here)
for 2 iterations,
but don't trust
this value in
general.

$J_{in} = 2$
 $J_{out} = 5$
is sufficient

Fig. 6. Averaged BER of the B-CG. Block size is $K_B = K$ where $K = 24, N = 32$, and $E_b/N_0 = 12$ dB. We consider one, two, three, and four inner iterations and the point GS is included for reference.

- The delay appears to be an inherent N symbol times to accumulate the measurements, plus computation time. Actually, it's much better than that, as shown in the next section.
- Note that the approach accommodates long codes.