

am230hw3

jxue16

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1 P1

If $[p_0, p_1 \dots p_n]$ is linear dependent, then there is $\alpha_0 p_0 + \dots + \alpha_n p_n = 0$.
for $i=0,1,\dots,n$

$$0 = p_i^T A(\alpha_0 p_0 + \dots + \alpha_n p_n) = \alpha_i p_i^T A p_i$$

$$\alpha_i p_i^T A p_i \neq 0, \text{ so } \alpha_i = 0$$

It is controversial, so $[p_0, p_1 \dots p_n]$ is linear independent

2 P2

$$\beta_{k+1} = \frac{r_{k+1}^T A p_k}{p_k^T A p_k}$$

$$r_{k+1} - r_k = \alpha_k A p_k$$

$$\therefore r_{k+1}^T A p_k = r_{k+1}^T (r_{k+1} - r_k) / \alpha_k = r_{k+1}^T r_{k+1} / \alpha_k$$

$$p_k = -r_k + \beta_{k+1} p_{k-1}$$

$$r_k^T p_k = -r_k^T r_k$$

$$p_k^T A p_k = -p_k^T r_k$$

$$p_k^T A p_k = -p_k^T r_k / \alpha_k = r_k^T r_k / \alpha_k$$

$$\therefore \beta_{k+1} = \frac{r_{k+1}^T r_{k+1}}{r_k^T r_k}$$

3 P3

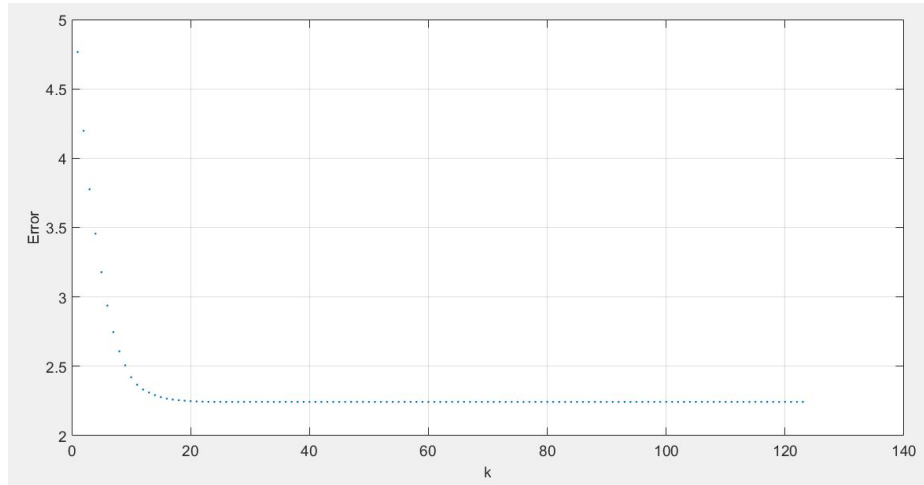


Figure 1

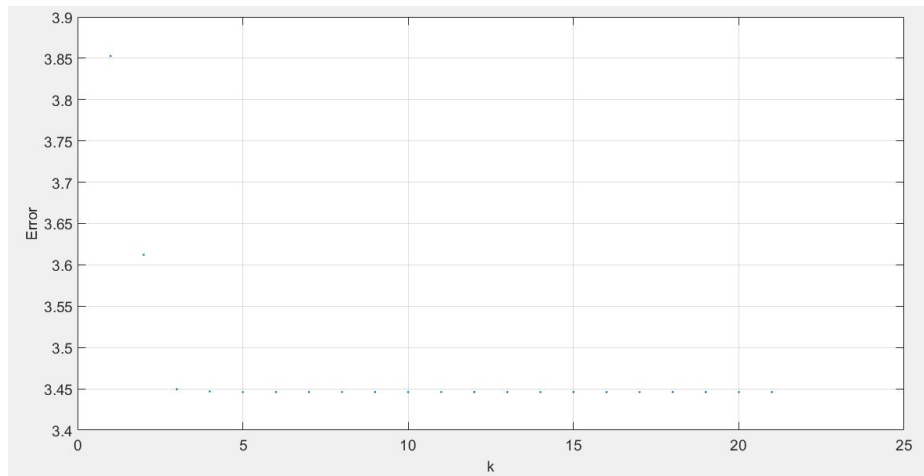


Figure 2

choose $n=100$

Figure 1 shows that without clustered eigenvalues, the convergence will be uniform. Figure 2 shows that the error of matrix A with two clustered eigenvalues will drop sharply after the some iterations.

4 P4

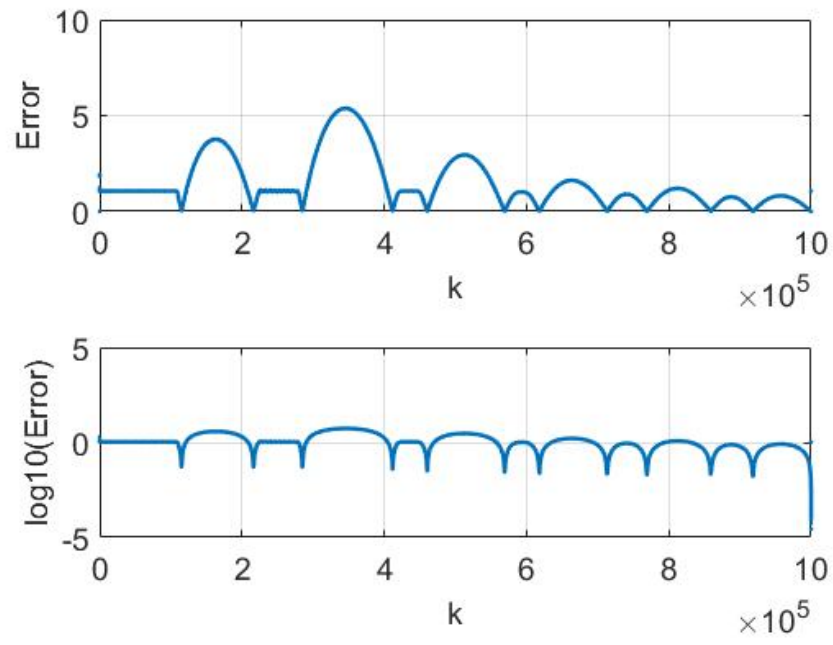


Figure 3: FR

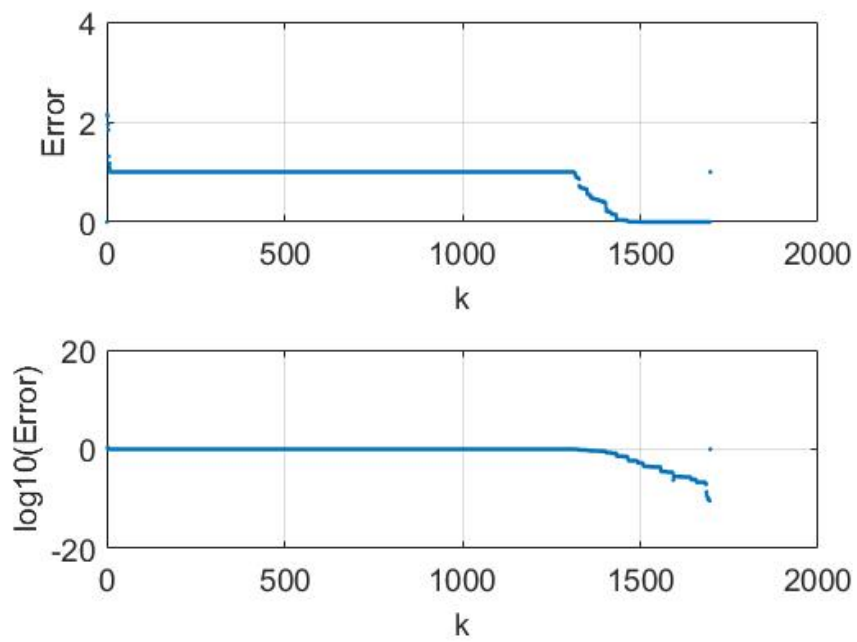


Figure 4: FR with restart

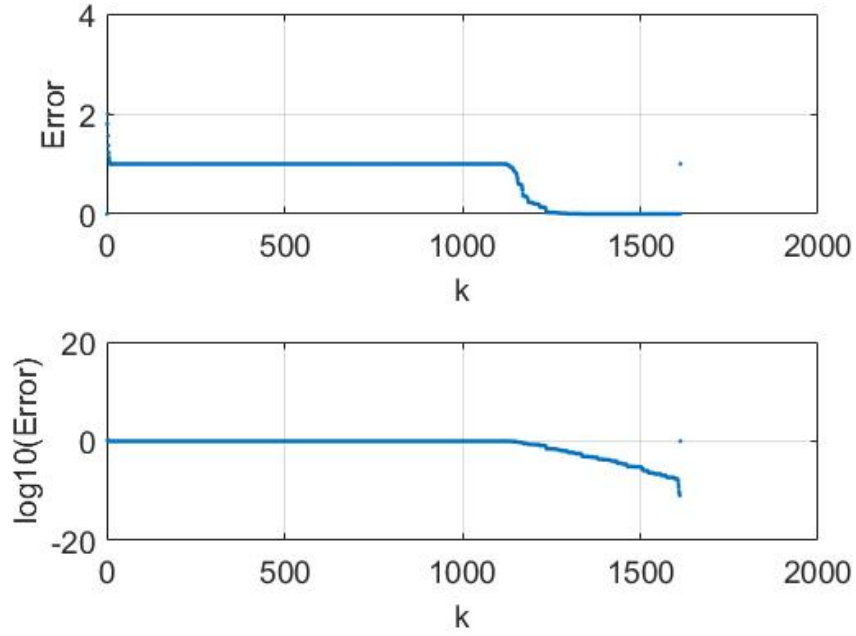


Figure 5: PR

Figure 1 shows that FR method doesn't converge well, it costs long time and the error plot oscillates.

Figure 2 shows that FR method can work well with restart. After 1600 iterations x can converge to $x = [1, 1, 1 \dots 1]^T$

Figure 3 shows that PR method can also work well without restart. So it is more robust and efficient than traditional FR method.