

Conjugate direction methods:

- a set of vectors, $p_0 \dots p_n$ is conjugate for A positive definite if $\checkmark \checkmark$

$$p_i' A p_j = 0 \quad i \neq j$$

- Assume we wish to min

$$\frac{x' A x}{2} - b' x$$

- useful because:

a) solution to $Ax = b$
for A p.d.

b) $\min_x \|Ax - b\|^2$ is like this

- now write

$$x = \alpha_0 p_0 + \alpha_1 p_1 + \dots$$

We then have

$$x'Ax + v'x = \alpha_0^2 p_0'Ap_0 + \alpha_1^2 p_1'Ap_1 + \dots + \alpha_0 v_0'p_0 + \alpha_1 v_1'p_1 + \dots$$

- we have no $p_i'Ap_j$ terms, the basis elements do not interact

- so we could minimize each α_i independently!

- conjugate dirns come from Gram-Schmidt

$$w_0 = e_0$$

$$w_1 = e_1 - \frac{(e_0'Aw_1)}{(e_0'Aw_0)} w_0$$

$$u_1 = \frac{w_1}{(w_1'Aw_1)^{1/2}}$$

$$w_2 = e_2 - (w_0'Aw_2)w_0 - (w_1'Aw_2)w_1$$

$$u_2 = \frac{w_2}{(w_2'Aw_2)^{1/2}}$$

Conjugate Direction in incremental form

Start with x_0, p_0

$$x_1 = x_0 + \alpha_0 p_0$$

now min wrt α_0

to get

$$\frac{(Ax_0 - b)' p_0}{p_0' A p_0} = \alpha_0$$

write

$$r_k = (Ax_k - b)$$

and get

$$x_{k+1} = x_k + \alpha_k p_k$$

$$\alpha_k = \frac{r_k' p_k}{p_k' A p_k}$$

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

properties of conjugate direction

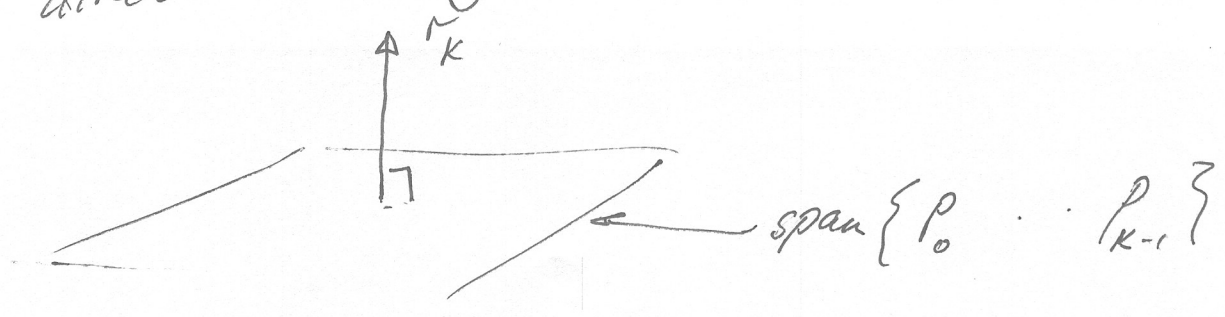
(i)

$$r_k^T \cdot p_i = 0 \quad \text{for } i < k$$

Proof: easy induction, p 16 a.

Significance:

- r_k is \perp to space spanned by p 's we have seen
- this means that ~~we have minimized~~ r_k might be a good source of a \perp direction conj to all other p 's.



props of C.D.

(13c)

(2) at k , we have minimized the quadratic form in $\text{span} \{p_0 \dots p_{k-1}\}$.

→ Obvious from Defn of conj. dirs., but v. important.

→ options: - go some steps, then keep x_k ; this has reduced the Q form value; if we are minimizing, this gives as a descent dir

→ What happens if A is not PD?

⇒ there is some $k+1$ so that the x_k is a max in that dir (or where the q form is constant!)

i.e. we choose

$$x_{k+1} = x_k + \alpha_k P_k$$

$$(x_k + \alpha_k P_k)^T A (x_k + \alpha_k P_k) = \alpha_k^2 (P_k^T A P_k) + 2 \alpha_k (x_k^T A P_k) + (x_k^T A x_k)$$

Now if $P_k^T A P_k > 0$, then this has a min otherwise, doesn't.

Strategy: go till we find a bad α_k ; keep the x_k , which is a better point than x_0

Minimizing: regard x_k as offset so current point.

No help in current form
because we need to know all
the p_i 's for a new direction.

But special choices of initial dir,
etc can make this unnecessary

Conjugate Gradient:

- want to solve

$$Ax = b$$

↑
symmetric, positive definite

equiv:

$$x = \operatorname{argmin}_x \frac{x'Ax}{2} - b'x$$

We will do this by generating special conj directions.

Start with

$$x_0$$

estimate of soln

$$r_0 = Ax_0 - b$$

residual

$$p_0 = -r_0$$

first dir

we will update by :

$$x_{k+1} = x_k + \alpha_k p_k$$

and get α_k by minimizing along p_k alone (OK, because p_k 's will be conjugate)

Again
this gives

$$\frac{1}{2} \left[(x_k + \alpha_k p_k)' A (x_k + \alpha_k p_k) \right] - b_k (x_k + \alpha_k p_k)$$

min is at:

$$\frac{-(Ax_k - b)' p_k}{p_k' A p_k}$$

write

$$r_k = Ax_k - b$$

$$\text{so } \alpha_k = \frac{-r_k' p_k}{p_k' A p_k}$$

Conjugate gradient

(15a)

(simple form)

Start

$$x_0, r_0 = Ax_0 - b, p_0 = -r_0$$

Step :

$$x_{k+1} = x_k + \alpha_k p_k$$

$$\alpha_k = \frac{-r_k' p_k}{p_k' A p_k}$$

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

This is
conj. direction
~~that~~
method

The clever
bit

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

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

from

$$\left[\leftarrow p_{k+1}' A p_k = 0 \right]$$

Properties of conjugate gradient

(166)

(1)

$$r_k' p_i = 0 \quad \text{for } i < k$$

assuming P conj.

Proof:

easy induction; notice

$$r_1' p_0 = 0 \quad (\text{by defn of } \alpha_0)$$
$$= r_0' r_0 + \alpha_0 p_0' A p_0$$

$$r_2' p_1 = 0 \quad ; \quad r_{k+1}' p_k = 0$$

then $r_2' p_0 = (r_1 + \alpha_1 A p_1)' p_0 = 0$ (because p are conj.)

(2)

$$r_k^T r_i = 0 \quad \text{for } i < k$$

Proof:

see 5.3 attached

(3)

$$\text{span} \{ r_0 \dots r_k \} = \text{span} \{ r_0, A r_0, \dots, A^k r_0 \}$$

(4)

$$\text{span} \{ p_0 \dots p_k \} = \text{span} \{ r_0, A r_0, \dots, A^k r_0 \}$$

(5)

$$p_k^T A p_i = 0$$

2A
Conjugate gradient.

Cleaner form:

• By properties, we have

$$\alpha_{k+1} = \frac{\Gamma_k' \Gamma_k}{P_k' A P_k}$$

• Now $\alpha_k A P_k = \Gamma_{k+1} - \Gamma_k$

$$\text{So } \beta_{k+1} = \Gamma_{k+1}' \left(\frac{\Gamma_{k+1} - \Gamma_k}{\alpha_k} \right) \cdot \frac{1}{P_k' A P_k}$$

$$= \frac{\Gamma_{k+1}' (\Gamma_{k+1} - \Gamma_k)}{\Gamma_k' \Gamma_k}$$

$$= \frac{\Gamma_{k+1}' \Gamma_{k+1}}{\Gamma_k' \Gamma_k} \quad (\text{By properties})$$

C.G. , clean form :

Start: x_0 , $r_0 = Ax_0 - b$, $P_0 = -r$

Step:

$$x_{k+1} = x_k + \alpha_k P_k$$

$$\alpha_k = \frac{r_k^T M_k}{P_k^T A P_k}$$

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

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

$$P_{k+1} = -r_{k+1} + \beta_{k+1} P_k$$

C.G. , properties.

if A is $n \times n$, n steps give \min .

in practice, can get to a good point rather fast; we might successfully n -steps; we might have a small residual early, too.

if A has r distinct eigenvalues, takes only r steps. !

"if about r eigenvalues form r "clusters", get v. good estimate. r steps are enough to

Potential applications

- Newton, Quasi Newton methods to solve linear alg.
- ~~Newton, where~~

- Large scale Newton, where we get

H_p with a Directional Derivative

Direct min. (Polak-Ribiere, etc).

- Keep doing CG until

x_k is a min

- Some issues to handle here

of \hat{x} . Because of the relation (5.9), however, the i th coordinate direction in \hat{x} -space corresponds to the direction p_i in x -space. Hence, the coordinate search strategy applied to $\hat{\phi}$ is equivalent to the conjugate direction algorithm (5.6), (5.7). We conclude, as in Theorem 5.1, that the conjugate direction algorithm terminates in at most n steps.

Returning to Figure 5.1, we note another interesting property: When the Hessian matrix is diagonal, each coordinate minimization correctly determines one of the components of the solution x^* . In other words, after k one-dimensional minimizations, the quadratic has been minimized on the subspace spanned by e_1, e_2, \dots, e_k . The following theorem proves this important result for the general case in which the Hessian of the quadratic is not necessarily diagonal. (Here and later, we use the notation $\text{span}\{p_0, p_1, \dots, p_k\}$ to denote the set of all linear combinations of the vectors p_0, p_1, \dots, p_k .) In proving the result we will make use of the following expression, which is easily verified from the relations (5.4) and (5.6):

$$r_{k+1} = r_k + \alpha_k A p_k. \tag{5.10}$$

Theorem 5.2 (Expanding Subspace Minimization).

Let $x_0 \in \mathbb{R}^n$ be any starting point and suppose that the sequence $\{x_k\}$ is generated by the conjugate direction algorithm (5.6), (5.7). Then

$$r_k^T p_i = 0, \quad \text{for } i = 0, 1, \dots, k-1, \tag{5.11}$$

and x_k is the minimizer of $\phi(x) = \frac{1}{2}x^T A x - b^T x$ over the set

$$\{x \mid x = x_0 + \text{span}\{p_0, p_1, \dots, p_{k-1}\}\}. \tag{5.12}$$

PROOF. We begin by showing that a point \tilde{x} minimizes ϕ over the set (5.12) if and only if $r(\tilde{x})^T p_i = 0$, for each $i = 0, 1, \dots, k-1$. Let us define $h(\sigma) = \phi(x_0 + \sigma_0 p_0 + \dots + \sigma_{k-1} p_{k-1})$, where $\sigma = (\sigma_0, \sigma_1, \dots, \sigma_{k-1})^T$. Since $h(\sigma)$ is a strictly convex quadratic, it has a unique minimizer σ^* that satisfies

$$\frac{\partial h(\sigma^*)}{\partial \sigma_i} = 0, \quad i = 0, 1, \dots, k-1.$$

By the chain rule, this equation implies that

$$\nabla\phi(x_0 + \sigma_0^* p_0 + \dots + \sigma_{k-1}^* p_{k-1})^T p_i = 0, \quad i = 0, 1, \dots, k-1.$$

By recalling the definition (5.3), we have for the minimizer $\tilde{x} = x_0 + \sigma_0^* p_0 + \sigma_1^* p_1 + \dots + \sigma_{k-1}^* p_{k-1}$ on the set (5.12) that $r(\tilde{x})^T p_i = 0$, as claimed.

We now use induction to show that x_k satisfies (5.11). For the case $k = 1$, we have from the fact that $x_1 = x_0 + \alpha_0 p_0$ minimizes ϕ along p_0 that $r_1^T p_0 = 0$. Let us now make

the induction hypothesis, namely, that $r_{k-1}^T p_i = 0$ for $i = 0, 1, \dots, k-2$. By (5.10), we have

$$r_k = r_{k-1} + \alpha_{k-1} A p_{k-1},$$

so that

$$p_{k-1}^T r_k = p_{k-1}^T r_{k-1} + \alpha_{k-1} p_{k-1}^T A p_{k-1} = 0,$$

by the definition (5.7) of α_{k-1} . Meanwhile, for the other vectors p_i , $i = 0, 1, \dots, k-2$, we have

$$p_i^T r_k = p_i^T r_{k-1} + \alpha_{k-1} p_i^T A p_{k-1} = 0,$$

where $p_i^T r_{k-1} = 0$ because of the induction hypothesis and $p_i^T A p_{k-1} = 0$ because of conjugacy of the vectors p_i . We have shown that $r_k^T p_i = 0$, for $i = 0, 1, \dots, k-1$, so the proof is complete. \square

The fact that the current residual r_k is orthogonal to all previous search directions, as expressed in (5.11), is a property that will be used extensively in this chapter.

The discussion so far has been general, in that it applies to a conjugate direction method (5.6), (5.7) based on *any* choice of the conjugate direction set $\{p_0, p_1, \dots, p_{n-1}\}$. There are many ways to choose the set of conjugate directions. For instance, the eigenvectors v_1, v_2, \dots, v_n of A are mutually orthogonal as well as conjugate with respect to A , so these could be used as the vectors $\{p_0, p_1, \dots, p_{n-1}\}$. For large-scale applications, however, computation of the complete set of eigenvectors requires an excessive amount of computation. An alternative approach is to modify the Gram-Schmidt orthogonalization process to produce a set of conjugate directions rather than a set of orthogonal directions. (This modification is easy to produce, since the properties of conjugacy and orthogonality are closely related in spirit.) However, the Gram-Schmidt approach is also expensive, since it requires us to store the entire direction set.

BASIC PROPERTIES OF THE CONJUGATE GRADIENT METHOD

The conjugate gradient method is a conjugate direction method with a very special property: In generating its set of conjugate vectors, it can compute a new vector p_k by using only the previous vector p_{k-1} . It does *not* need to know all the previous elements p_0, p_1, \dots, p_{k-2} of the conjugate set; p_k is automatically conjugate to these vectors. This remarkable property implies that the method requires little storage and computation.

In the conjugate gradient method, each direction p_k is chosen to be a linear combination of the negative residual $-r_k$ (which, by (5.3), is the steepest descent direction for the

function ϕ) and the previous direction p_{k-1} . We write

$$p_k = -r_k + \beta_k p_{k-1}, \quad (5.13)$$

where the scalar β_k is to be determined by the requirement that p_{k-1} and p_k must be conjugate with respect to A . By premultiplying (5.13) by $p_{k-1}^T A$ and imposing the condition $p_{k-1}^T A p_k = 0$, we find that

$$\beta_k = \frac{r_k^T A p_{k-1}}{p_{k-1}^T A p_{k-1}}.$$

We choose the first search direction p_0 to be the steepest descent direction at the initial point x_0 . As in the general conjugate direction method, we perform successive one-dimensional minimizations along each of the search directions. We have thus specified a complete algorithm, which we express formally as follows:

Algorithm 5.1 (CG–Preliminary Version).

Given x_0 ;
Set $r_0 \leftarrow Ax_0 - b$, $p_0 \leftarrow -r_0$, $k \leftarrow 0$;
while $r_k \neq 0$

$$\alpha_k \leftarrow -\frac{r_k^T p_k}{p_k^T A p_k}; \quad (5.14)$$

$$x_{k+1} \leftarrow x_k + \alpha_k p_k; \quad (5.14)$$

$$r_{k+1} \leftarrow Ax_{k+1} - b; \quad (5.14)$$

$$\beta_{k+1} \leftarrow \frac{r_{k+1}^T A p_k}{p_k^T A p_k}; \quad (5.14)$$

$$p_{k+1} \leftarrow -r_{k+1} + \beta_{k+1} p_k; \quad (5.14)$$

$$k \leftarrow k + 1; \quad (5.14)$$

end (while)

This version is useful for studying the essential properties of the conjugate gradient method, but we present a more efficient version later. We show first that the directions p_0, p_1, \dots, p_{n-1} are indeed conjugate, which by Theorem 5.1 implies termination in n steps. The theorem below establishes this property and two other important properties. First, the residuals r_i are mutually orthogonal. Second, each search direction p_k and residual r_k is contained in the *Krylov subspace of degree k for r_0* , defined as

$$\mathcal{K}(r_0; k) \stackrel{\text{def}}{=} \text{span}\{r_0, Ar_0, \dots, A^k r_0\}. \quad (5.15)$$

Theorem 5.3.

Suppose that the k th iterate generated by the conjugate gradient method is not the solution point x^* . The following four properties hold:

$$r_k^T r_i = 0, \quad \text{for } i = 0, 1, \dots, k-1, \quad (5.16)$$

$$\text{span}\{r_0, r_1, \dots, r_k\} = \text{span}\{r_0, Ar_0, \dots, A^k r_0\}, \quad (5.17)$$

$$\text{span}\{p_0, p_1, \dots, p_k\} = \text{span}\{r_0, Ar_0, \dots, A^k r_0\}, \quad (5.18)$$

$$p_k^T A p_i = 0, \quad \text{for } i = 0, 1, \dots, k-1. \quad (5.19)$$

Therefore, the sequence $\{x_k\}$ converges to x^* in at most n steps.

PROOF. The proof is by induction. The expressions (5.17) and (5.18) hold trivially for $k = 0$, while (5.19) holds by construction for $k = 1$. Assuming now that these three expressions are true for some k (the induction hypothesis), we show that they continue to hold for $k + 1$.

To prove (5.17), we show first that the set on the left-hand side is contained in the set on the right-hand side. Because of the induction hypothesis, we have from (5.17) and (5.18) that

$$r_k \in \text{span}\{r_0, Ar_0, \dots, A^k r_0\}, \quad p_k \in \text{span}\{r_0, Ar_0, \dots, A^k r_0\},$$

while by multiplying the second of these expressions by A , we obtain

$$A p_k \in \text{span}\{Ar_0, \dots, A^{k+1} r_0\}. \quad (5.20)$$

By applying (5.10), we find that

$$r_{k+1} \in \text{span}\{r_0, Ar_0, \dots, A^{k+1} r_0\}.$$

By combining this expression with the induction hypothesis for (5.17), we conclude that

$$\text{span}\{r_0, r_1, \dots, r_k, r_{k+1}\} \subset \text{span}\{r_0, Ar_0, \dots, A^{k+1} r_0\}.$$

To prove that the reverse inclusion holds as well, we use the induction hypothesis on (5.18) to deduce that

$$A^{k+1} r_0 = A(A^k r_0) \in \text{span}\{A p_0, A p_1, \dots, A p_k\}.$$

Since by (5.10) we have $A p_i = (r_{i+1} - r_i)/\alpha_i$ for $i = 0, 1, \dots, k$, it follows that

$$A^{k+1} r_0 \in \text{span}\{r_0, r_1, \dots, r_{k+1}\}.$$

By combining this expression with the induction hypothesis for (5.17), we find that

$$\text{span}\{r_0, Ar_0, \dots, A^{k+1}r_0\} \subset \text{span}\{r_0, r_1, \dots, r_k, r_{k+1}\}.$$

Therefore, the relation (5.17) continues to hold when k is replaced by $k + 1$, as claimed.

We show that (5.18) continues to hold when k is replaced by $k + 1$ by the following argument:

$$\begin{aligned} & \text{span}\{p_0, p_1, \dots, p_k, p_{k+1}\} \\ &= \text{span}\{p_0, p_1, \dots, p_k, r_{k+1}\} && \text{by (5.14e)} \\ &= \text{span}\{r_0, Ar_0, \dots, A^k r_0, r_{k+1}\} && \text{by induction hypothesis for (5.18)} \\ &= \text{span}\{r_0, r_1, \dots, r_k, r_{k+1}\} && \text{by (5.17)} \\ &= \text{span}\{r_0, Ar_0, \dots, A^{k+1}r_0\} && \text{by (5.17) for } k + 1. \end{aligned}$$

Next, we prove the conjugacy condition (5.19) with k replaced by $k + 1$. By multiplying (5.14e) by Ap_i , $i = 0, 1, \dots, k$, we obtain

$$p_{k+1}^T Ap_i = -r_{k+1}^T Ap_i + \beta_{k+1} p_k^T Ap_i. \quad (5.21)$$

By the definition (5.14d) of β_k , the right-hand-side of (5.21) vanishes when $i = k$. For $i \leq k - 1$ we need to collect a number of observations. Note first that our induction hypothesis for (5.19) implies that the directions p_0, p_1, \dots, p_k are conjugate, so we can apply Theorem 5.2 to deduce that

$$r_{k+1}^T p_i = 0, \quad \text{for } i = 0, 1, \dots, k. \quad (5.22)$$

Second, by repeatedly applying (5.18), we find that for $i = 0, 1, \dots, k - 1$, the following inclusion holds:

$$\begin{aligned} Ap_i &\in A \text{span}\{r_0, Ar_0, \dots, A^i r_0\} = \text{span}\{Ar_0, A^2 r_0, \dots, A^{i+1} r_0\} \\ &\subset \text{span}\{p_0, p_1, \dots, p_{i+1}\}. \end{aligned} \quad (5.23)$$

By combining (5.22) and (5.23), we deduce that

$$r_{k+1}^T Ap_i = 0, \quad \text{for } i = 0, 1, \dots, k - 1,$$

so the first term in the right-hand-side of (5.21) vanishes for $i = 0, 1, \dots, k - 1$. Because of the induction hypothesis for (5.19), the second term vanishes as well, and we

conclude that $p_{k+1}^T A p_i = 0, i = 0, 1, \dots, k$. Hence, the induction argument holds for (5.19) also.

It follows that the direction set generated by the conjugate gradient method is indeed a conjugate direction set, so Theorem 5.1 tells us that the algorithm terminates in at most n iterations.

Finally, we prove (5.16) by a noninductive argument. Because the direction set is conjugate, we have from (5.11) that $r_k^T p_i = 0$ for all $i = 0, 1, \dots, k-1$ and any $k = 1, 2, \dots, n-1$. By rearranging (5.14e), we find that

$$p_i = -r_i + \beta_i p_{i-1},$$

so that $r_i \in \text{span}\{p_i, p_{i-1}\}$ for all $i = 1, \dots, k-1$. We conclude that $r_k^T r_i = 0$ for all $i = 1, \dots, k-1$. To complete the proof, we note that $r_k^T r_0 = -r_k^T p_0 = 0$, by definition of p_0 in Algorithm 5.1 and by (5.11). \square

The proof of this theorem relies on the fact that the first direction p_0 is the steepest descent direction $-r_0$; in fact, the result does not hold for other choices of p_0 . Since the gradients r_k are mutually orthogonal, the term "conjugate gradient method" is actually a misnomer. It is the search directions, not the gradients, that are conjugate with respect to A .

A PRACTICAL FORM OF THE CONJUGATE GRADIENT METHOD

We can derive a slightly more economical form of the conjugate gradient method by using the results of Theorems 5.2 and 5.3. First, we can use (5.14e) and (5.11) to replace the formula (5.14a) for α_k by

$$\alpha_k = \frac{r_k^T r_k}{p_k^T A p_k}.$$

Second, we have from (5.10) that $\alpha_k A p_k = r_{k+1} - r_k$, so by applying (5.14e) and (5.11) once again we can simplify the formula for β_{k+1} to

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

By using these formulae together with (5.10), we obtain the following standard form of the conjugate gradient method.

Algorithm 5.2 (CG).Given x_0 ;Set $r_0 \leftarrow Ax_0 - b$, $p_0 \leftarrow -r_0$, $k \leftarrow 0$;**while** $r_k \neq 0$

$$\alpha_k \leftarrow \frac{r_k^T r_k}{p_k^T A p_k}; \quad (5.24a)$$

$$x_{k+1} \leftarrow x_k + \alpha_k p_k; \quad (5.24b)$$

$$r_{k+1} \leftarrow r_k + \alpha_k A p_k; \quad (5.24c)$$

$$\beta_{k+1} \leftarrow \frac{r_{k+1}^T r_{k+1}}{r_k^T r_k}; \quad (5.24d)$$

$$p_{k+1} \leftarrow -r_{k+1} + \beta_{k+1} p_k; \quad (5.24e)$$

$$k \leftarrow k + 1; \quad (5.24f)$$

end (while)

At any given point in Algorithm 5.2 we never need to know the vectors x , r , and p for more than the last two iterations. Accordingly, implementations of this algorithm overwrite old values of these vectors to save on storage. The major computational tasks to be performed at each step are computation of the matrix-vector product $A p_k$, calculation of the inner products $p_k^T (A p_k)$ and $r_{k+1}^T r_{k+1}$, and calculation of three vector sums. The inner product and vector sum operations can be performed in a small multiple of n floating-point operations, while the cost of the matrix-vector product is, of course, dependent on the problem. The CG method is recommended only for large problems; otherwise, Gaussian elimination or other factorization algorithms such as the singular value decomposition are to be preferred, since they are less sensitive to rounding errors. For large problems, the CG method has the advantage that it does not alter the coefficient matrix and (in contrast to factorization techniques) does not produce fill in the arrays holding the matrix. Another key property is that the CG method sometimes approaches the solution quickly, as we discuss next.

RATE OF CONVERGENCE

We have seen that in exact arithmetic the conjugate gradient method will terminate at the solution in at most n iterations. What is more remarkable is that when the distribution of the eigenvalues of A has certain favorable features, the algorithm will identify the solution in many fewer than n iterations. To explain this property, we begin by viewing the expanding subspace minimization property proved in Theorem 5.2 in a slightly different way, using it to show that Algorithm 5.2 is optimal in a certain important sense.