

# Conjugate gradient method

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Optimization methods. MIPT

## Strongly convex quadratics

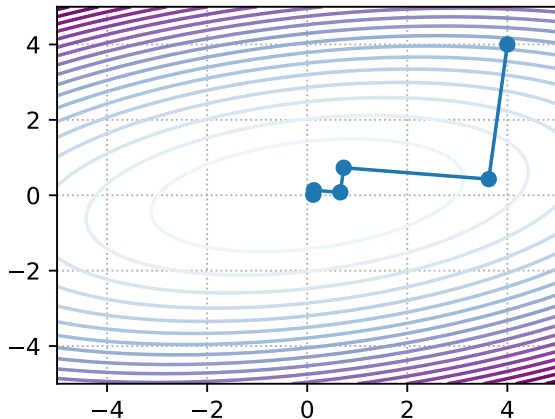
Consider the following quadratic optimization problem:

$$\min_{x \in \mathbb{R}^d} f(x) = \min_{x \in \mathbb{R}^d} \frac{1}{2} x^\top A x - b^\top x + c, \text{ where } A \in \mathbb{S}_{++}^d.$$

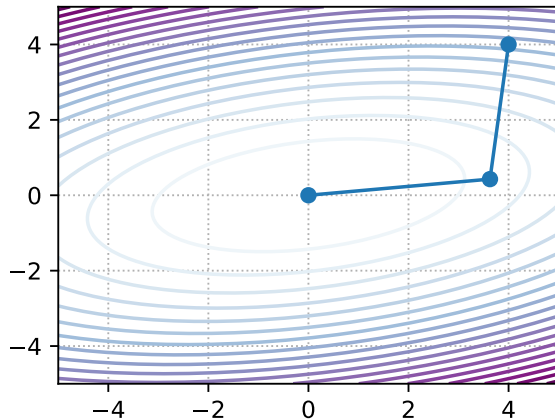
Optimality conditions

$$Ax^* = b$$

Steepest Descent



Conjugate Gradient



## Exact line search aka steepest descent

$$\alpha_k = \arg \min_{\alpha \in \mathbb{R}^+} f(x_{k+1}) = \arg \min_{\alpha \in \mathbb{R}^+} f(x_k - \alpha \nabla f(x_k))$$

More theoretical than practical approach. It also allows you to analyze the convergence, but often exact line search can be difficult if the function calculation takes too long or costs a lot. An interesting theoretical property of this method is that each following iteration is orthogonal to the previous one:

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Optimality conditions:

$$\nabla f(x_k)^T \nabla f(x_{k+1}) = 0$$

🔥 Optimal value for quadratics

$$\nabla f(x_k)^T A(x_k - \alpha \nabla f(x_k)) - \nabla f(x_k)^T b = 0 \quad \alpha_k = \frac{\nabla f(x_k)^T \nabla f(x_k)}{\nabla f(x_k)^T A \nabla f(x_k)}$$



Figure 1: Steepest Descent

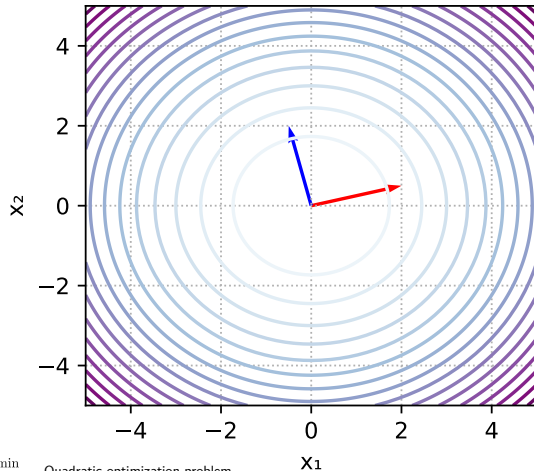
Open In Colab

## Conjugate directions. $A$ -orthogonality.

$v_1$  and  $v_2$  are orthogonal

$$v_1^T v_2 = 0.00$$

$$v_1^T A v_2 = 1.19$$



$\hat{v}_1$  and  $\hat{v}_2$  are  $A$ -orthogonal

$$\hat{v}_1^T \hat{v}_2 = -0.80$$

$$\hat{v}_1^T A \hat{v}_2 = -0.00$$



## Conjugate directions. $A$ -orthogonality.

Suppose, we have two coordinate systems and some quadratic function  $f(x) = \frac{1}{2}x^T I x$  looks just like on the left part of Figure 2, while in another coordinates it looks like  $f(\hat{x}) = \frac{1}{2}\hat{x}^T A \hat{x}$ , where  $A \in \mathbb{S}_{++}^d$ .

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Since  $A = Q\Lambda Q^T$ :

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### $A$ -orthogonal vectors

Vectors  $x \in \mathbb{R}^d$  and  $y \in \mathbb{R}^d$  are called  $A$ -orthogonal (or  $A$ -conjugate) if

$$x^T A y = 0 \quad \Leftrightarrow \quad x \perp_A y$$

When  $A = I$ ,  $A$ -orthogonality becomes orthogonality.

# Gram–Schmidt process

# Idea of the method of conjugate directions

Thus, we formulate an algorithm:

1. Let  $k = 0$  and  $x_k = x_0$ , count  $d_k = d_0 = -\nabla f(x_0)$ .



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$$\alpha_k = -\frac{d_k^\top (Ax_k - b)}{d_k^\top Ad_k}$$

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4. update the direction:  $d_{k+1} = -\nabla f(x_{k+1}) + \beta_k d_k$ , where  $\beta_k$  is calculated by the formula:

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5. Repeat steps 2-4 until  $n$  directions are built, where  $n$  is the dimension of space (dimension of  $x$ ).

## Method of Conjugate Directions

If a set of vectors  $d_1, \dots, d_k$  - are  $A$ -conjugate (each pair of vectors is  $A$ -conjugate), these vectors are linearly independent.  $A \in \mathbb{S}_{++}^n$ .

### Proof

We'll show, that if  $\sum_{i=1}^k \alpha_i d_i = 0$ , than all coefficients should be equal to zero:

$$\begin{aligned} 0 &= \sum_{i=1}^n \alpha_i d_i \\ &= d_j^\top A \left( \sum_{i=1}^n \alpha_i d_i \right) \\ &= \sum_{i=1}^n \alpha_i d_j^\top A d_i \\ &= \alpha_j d_j^\top A d_j + 0 + \dots + 0 \end{aligned}$$

Thus,  $\alpha_j = 0$ , for all other indices one have perform the same process

# Conjugate Gradients

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# Conjugate gradient method



Conjugate Gradient = Conjugate Directions  
+ Residuals as starting vectors for Gram–Schmidt

$$\mathbf{r}_0 := \mathbf{b} - \mathbf{A}\mathbf{x}_0$$

if  $\mathbf{r}_0$  is sufficiently small, then return  $\mathbf{x}_0$  as the result

$$\mathbf{d}_0 := \mathbf{r}_0$$

$$k := 0$$

repeat

$$\alpha_k := \frac{\mathbf{r}_k^\top \mathbf{r}_k}{\mathbf{d}_k^\top \mathbf{A} \mathbf{d}_k}$$

$$\mathbf{x}_{k+1} := \mathbf{x}_k + \alpha_k \mathbf{d}_k$$

$$\mathbf{r}_{k+1} := \mathbf{r}_k - \alpha_k \mathbf{A} \mathbf{d}_k$$

if  $\mathbf{r}_{k+1}$  is sufficiently small, then exit loop

$$\beta_k := \frac{\mathbf{r}_{k+1}^\top \mathbf{r}_{k+1}}{\mathbf{r}_k^\top \mathbf{r}_k}$$

$$\mathbf{d}_{k+1} := \mathbf{r}_{k+1} + \beta_k \mathbf{d}_k$$

$$k := k + 1$$

end repeat

return  $\mathbf{x}_{k+1}$  as the result

# Convergence

**Theorem 1.** If matrix  $A$  has only  $r$  different eigenvalues, then the conjugate gradient method converges in  $r$  iterations.

**Theorem 2.** The following convergence bound holds

$$\|x_k - x^*\|_A \leq 2 \left( \frac{\sqrt{\kappa(A)} - 1}{\sqrt{\kappa(A)} + 1} \right)^k \|x_0 - x^*\|_A,$$

where  $\|x\|_A^2 = x^\top A x$  and  $\kappa(A) = \frac{\lambda_1(A)}{\lambda_n(A)}$  is the conditioning number of matrix  $A$ ,  $\lambda_1(A) \geq \dots \geq \lambda_n(A)$  are the eigenvalues of matrix  $A$

**Note:** compare the coefficient of the geometric progression with its analog in gradient descent.

## Non-linear conjugate gradient method

In case we do not have an analytic expression for a function or its gradient, we will most likely not be able to solve the one-dimensional minimization problem analytically. Therefore, step 2 of the algorithm is replaced by the usual line search procedure. But there is the following mathematical trick for the fourth point:

For two iterations, it is fair:

$$x_{k+1} - x_k = cd_k,$$

where  $c$  is some kind of constant. Then for the quadratic case, we have:

$$\nabla f(x_{k+1}) - \nabla f(x_k) = (Ax_{k+1} - b) - (Ax_k - b) = A(x_{k+1} - x_k) = cAd_k$$

Expressing from this equation the work  $Ad_k = \frac{1}{c} (\nabla f(x_{k+1}) - \nabla f(x_k))$ , we get rid of the “knowledge” of the function in step definition  $\beta_k$ , then point 4 will be rewritten as:

$$\beta_k = \frac{\nabla f(x_{k+1})^\top (\nabla f(x_{k+1}) - \nabla f(x_k))}{d_k^\top (\nabla f(x_{k+1}) - \nabla f(x_k))}.$$

This method is called the Polack - Ribier method.

# Preconditioned conjugate gradient method