# **Introduction to Machine Learning**

# **ML-Basics Optimization**





#### **Learning goals**

- $\bullet$  Understand how the risk function is optimized to learn the optimal parameters of a model
- Understand the idea of gradient descent as a basic risk optimizer

# **LEARNING AS PARAMETER OPTIMIZATION**

- We have seen, we can operationalize the search for a model *f* that matches training data best, by looking for its parametrization  $\theta \in \Theta$  with lowest empirical risk  $\mathcal{R}_{emp}(\theta)$ .
- Therefore, we usually traverse the error surface downwards; often by local search from a starting point to its minimum.



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### **LEARNING AS PARAMETER OPTIMIZATION / 2**

The ERM optimization problem is:

$$
\hat{\theta} = \argmin_{\theta \in \Theta} \mathcal{R}_{\textsf{emp}}(\theta).
$$

For a **(global) minimum**  $\hat{\theta}$  it obviously holds that

$$
\forall \theta \in \Theta: \quad \mathcal{R}_{\text{emp}}(\hat{\theta}) \leq \mathcal{R}_{\text{emp}}(\theta).
$$

This does not imply that  $\hat{\theta}$  is unique.

Which kind of numerical technique is reasonable for this problem strongly depends on model and parameter structure (continuous params? uni-modal  $\mathcal{R}_{\text{emo}}(\theta)$ ?). Here, we will only discuss very simple scenarios.

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#### **LOCAL MINIMA**

If  $\mathcal{R}_{\text{emo}}$  is continuous in  $\theta$  we can define a **local minimum**  $\hat{\theta}$ :

$$
\exists \epsilon > 0 \; \forall \pmb{\theta} \text{ with } \left\|\hat{\pmb{\theta}} - \pmb{\theta}\right\| < \epsilon: \quad \mathcal{R}_{\sf emp}(\hat{\pmb{\theta}}) \leq \mathcal{R}_{\sf emp}(\pmb{\theta}).
$$

Clearly every global minimum is also a local minimum. Finding a local minimum is easier than finding a global minimum.





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# **LOCAL MINIMA AND STATIONARY POINTS**

If  $\mathcal{R}_{\text{emo}}$  is continuously differentiable in  $\theta$  then a **sufficient condition** for a local minimum is that  $\hat{\theta}$  is **stationary** with 0 gradient, so no local improvement is possible:

$$
\frac{\partial}{\partial\boldsymbol{\theta}}\mathcal{R}_{\mathsf{emp}}(\boldsymbol{\hat{\theta}})=0
$$

and the Hessian  $\frac{\partial^2}{\partial\bm{\theta}^2}\mathcal{R}_{\sf emp}(\hat{\bm{\theta}})$  is positive definite. While the neg. gradient points into the direction of fastest local decrease, the Hessian measures local curvature of  $\mathcal{R}_{\text{emo}}$ .



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## **LEAST SQUARES ESTIMATOR**

Now, for given features  $\mathbf{X} \in \mathbb{R}^{n \times p}$  and target  $\mathbf{y} \in \mathbb{R}^{n}$ , we want to find the best linear model regarding the squared error loss, i.e.,

$$
\mathcal{R}_{\text{emp}}(\boldsymbol{\theta}) = \|\mathbf{X}\boldsymbol{\theta} - \mathbf{y}\|_2^2 = \sum_{i=1}^n (\boldsymbol{\theta}^\top \mathbf{x}^{(i)} - y^{(i)})^2.
$$

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With the sufficient condition for continously differentiable functions it can be shown that the **least squares estimator**

$$
\hat{\theta} = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top \mathbf{y}.
$$

is a local minimum of  $\mathcal{R}_{\text{emp}}$ . If **X** is full-rank,  $\mathcal{R}_{\text{emp}}$  is strictly convex and there is only one local minimum - which is also global.

**Note:** Often such analytical solutions in ML are not possible, and we rather have to use iterative numerical optimization.

#### **GRADIENT DESCENT**

The simple idea of GD is to iteratively go from the current candidate  $\theta^{[t]}$ in the direction of the negative gradient, i.e., the direction of the steepest descent, with learning rate  $\alpha$  to the next  $\boldsymbol{\theta}^{[t+1]}.$ 

$$
\boldsymbol{\theta}^{[t+1]} = \boldsymbol{\theta}^{[t]} - \alpha \frac{\partial}{\partial \boldsymbol{\theta}} \mathcal{R}_{\text{emp}}(\boldsymbol{\theta}^{[t]}).
$$





We choose a random start  $\theta^{[0]}$  with risk  $\mathcal{R}_{\mathsf{emp}}(\boldsymbol{\theta}^{[0]})=76.25.$ 

#### **GRADIENT DESCENT - EXAMPLE**



Now we follow in the direction of the negative gradient at  $\theta^{[0]}$ .

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We arrive at  $\theta^{[1]}$  with risk  $\mathcal{R}_{\mathsf{emp}}(\boldsymbol{\theta}^{[1]}) \approx$  42.73. We improved:  $\mathcal{R}_{\mathsf{emp}}(\boldsymbol{\theta}^{[1]}) < \mathcal{R}_{\mathsf{emp}}(\boldsymbol{\theta}^{[0]}).$ 

#### **GRADIENT DESCENT - EXAMPLE**



Again we follow in the direction of the negative gradient, but now at  $\boldsymbol{\theta}^{[1]}$ .

Now  $\theta^{[2]}$  has risk  $\mathcal{R}_{\mathsf{emp}}(\theta^{[2]})\approx$  25.08.

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#### **GRADIENT DESCENT - EXAMPLE**



We iterate this until some form of convergence or termination.

We arrive close to a stationary  $\hat{\theta}$  which is hopefully at least a local minimum.

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# **GRADIENT DESCENT - LEARNING RATE**

- $\bullet$  The negative gradient is a direction that looks locally promising to reduce  $\mathcal{R}_{\text{emo}}$ .
- $\bullet$  Hence it weights components higher in which  $\mathcal{R}_{\text{emo}}$  decreases more.
- However, the length of  $-\frac{\partial}{\partial \theta} \mathcal{R}_{\text{\rm emp}}$  measures only the local decrease rate, i.e., there are no guarantees that we will not go "too far".
- $\bullet$  We use a learning rate  $\alpha$  to scale the step length in each iteration. Too much can lead to overstepping and no converge, too low leads to slow convergence.
- Usually, a simple constant rate or rate-decrease mechanisms to enforce local convergence are used





#### **FURTHER TOPICS**

- GD is a so-called first-order method. Second-order methods use the Hessian to refine the search direction for faster convergence.
- There exist many improvements of GD, e.g., to smartly control the learn rate, to escape saddle points, to mimic second order behavior without computing the expensive Hessian.
- If the gradient of GD is not derived from the empirical risk of the whole data set, but instead from a randomly selected subset, we call this **stochastic gradient descent** (SGD). For large-scale problems this can lead to higher computational efficiency.