Chapter 1 Linear Regression

> **Introduction**
>
> ❏ *ERM* ❏ *Ridge Regression (L_2 regularization)*
>
> ❏ *Gradient Descent* ❏ *Lasso Regression (L_1 regularization)*

1.1 Basic Knowledge

Example 1.1 Linear Regression

Settings.

- Dataset: $D = \{(x_i, y_i)\}_{i=1}^n$, where $x_i \in \mathbb{R}^d$ and $y_i \in \mathbb{R}$. Here, y_i denotes the regression target, while x_i represents the input features used to predict y_i.
- Linear Model: $f(x) = w^\top x + b$, with weight $w \in \mathbb{R}^d$ and bias $b \in \mathbb{R}$.
- Linear Model: $f(x) = w^\top x + b$, with weight $w \in \mathbb{R}^d$ and bias $b \in \mathbb{R}$.

⚲ **Note** *This definition is equivalent to an inner product: $\hat{y} = w^\top x + b$. $\hat{y} = w^\top x + b$.*

> **Definition 1.1 (Learnable / Trainable Parameters)**
>
> *Learnable parameters are those that can be updated during the training process.* ♣

Quiz. How to determine whether a parameter is learnable? Quiz: How to determine w and b? Quiz: How to determine w and b?

Ans: **ERM** (Empirical Risk Minimization)

- Loss function. Squared Loss (SE) is commonly used during optimization. The training objective can be written as:

$$\underset{w,b}{\text{argmin}} \frac{1}{n} \sum_{i \in [n]} \left(y_i - (w^\top x_i + b)\right)^2 \underset{w,b}{\text{argmin}} \frac{1}{n} \sum_{i \in [n]} \left(y_i - (w^\top x_i + b)\right)^2 \tag{1.1}$$

The blue factor $1/n$ can be omitted in theoretical analysis, but is often kept in practice to stabilize the loss function during implementation.

Quiz: How to optimize the parameters?

Ans: **Gradient Descent** (as a traditional ML method). In the case of linear regression:

$$\frac{\partial \mathcal{L}}{\partial b} = -2 \sum_{i \in [n]} (y_i - w^\top x_i - b) \tag{1.2}$$

$$\frac{\partial \mathcal{L}}{\partial w} = -2 \sum_{i \in [n]} (y_i - w^\top x_i - b)x_i \tag{1.3}$$

⚲ **Note** *In the field of machine learning, the gradient of a scalar with respect to a vector is itself a vector (**not a covector**). This means:*

$$\frac{\partial \mathcal{L}}{\partial w} = \begin{pmatrix} \frac{\partial \mathcal{L}}{\partial w_1} \\ \frac{\partial \mathcal{L}}{\partial w_2} \\ \vdots \\ \frac{\partial \mathcal{L}}{\partial w_d} \end{pmatrix} = \left(\frac{\partial \mathcal{L}}{\partial w_1}, \frac{\partial \mathcal{L}}{\partial w_2}, \cdots, \frac{\partial \mathcal{L}}{\partial w_d} \right)^\top \tag{1.4}$$

See the definition of matix derivatives (assuming no special structure in the matrix) in Matrix Cookbook Chapter 2.

Note *Here are some commonly used derivative formulas:*

$$\frac{\partial x^\top x}{\partial x} = 2x \tag{1.5}$$

$$\frac{\partial a^\top x}{\partial x} = a, \quad \frac{\partial Ax}{\partial x} = A^\top \tag{1.6}$$

$$\frac{\partial x^\top Ax}{\partial x} = (A + A^\top)x \tag{1.7}$$

Remark Both sides of an equation must have the same dimension. This principle can be used as a consistency check.

We optimize the parameters by subtracting a scalar multiple of the gradient from the parameters, considering the physical meaning of the gradient: the direction of the steepest **increase**.

> **Definition 1.2 (Hyperparameter)**
>
> *A parameter that is fixed during optimization and specified before the training process.* ♣

That is:

$$w' = w - \alpha \frac{\partial \mathcal{L}}{\partial w}, \quad b' = b - \alpha \frac{\partial \mathcal{L}}{\partial b} \tag{1.8}$$

Optimization will stop when the norm of the parameter update becomes smaller than a given hyperparameter.

1.2 Closed-Form of Linear Regression

> **Proposition 1.1**
>
> *Linear Regression has **Closed-Form** solution.* ♡

Settings.
- Matrix $X_0 := (x_1, \cdots, x_n)^\top$;
- Matrix $X := (X_0, \mathbb{1}) \in \mathbb{R}^{n \times (d+1)}$;
- $y = (y_1, \cdots, y_n)^\top \in \mathbb{R}^n$;
- $\hat{w} = (w^\top, b)^\top \in \mathbb{R}^{d+1}$.

Then the loss function of $\hat{w}$ can be written as:

$$\mathcal{L}(\hat{w}) = (y - X\hat{w})^\top (y - X\hat{w}) = \|y - X\hat{w}\|_2^2 \tag{1.9}$$

Here, $\| \cdot \|_p$ denotes the *p-norm* of a vector.

Note *Vectors can sometimes be treated as scalars, since linearity ensures that the validity of a proposition can be extended to any finite dimension.*

Notice that the optimization stops when $\partial \mathcal{L}(\hat{w})/\partial \hat{w} = 0$. Under this condition, the parameters can be solved from the above constraint by following steps:

$$\frac{\partial \mathcal{L}(\hat{w})}{\partial \hat{w}} = -2X^\top (y - X\hat{w}) \tag{1.10}$$

Note *Both dimensional analysis and calculation using Leibniz's rule lead to the same result as the formula above:*

$$\mathcal{L}(\hat{w}) = y^\top y - 2y^\top X\hat{w} + \hat{w}^\top X^\top X\hat{w}$$

$$\partial_{\hat{w}} \mathcal{L}(\hat{w}) = -2X^\top y + 2X^\top X\hat{w}$$

$$= -2X^\top (y - X\hat{w})$$

Remark More matix formulas are available in Matrix Cookbook.

Thus, the target of the optimization satisfied:

$$X^\top y = X^\top X\hat{w} \tag{1.11}$$

That is:

$$\hat{w} = (X^\top X)^{-1} X^\top y \tag{1.12}$$

when $X^\top X$ invertible (non-singular / full-rank).

Example 1.2 When does $X^\top X$ not invertible?

Solution $X \in \mathbb{R}^{n \times (d+1)}$:

- $d + 1 > n$. *Brief Proof:* $\text{rank}(X^\top X) = \text{rank}(X) \leq \min(n, d+1) = n < d+1$.
- *X has repeated columns. Proof is trivial.*

When $X^\top X$ isn't invertible:

1. If $\text{rank}(X^\top X, X^\top y) > \text{rank}(X^\top X)$, $\hat{w}$ has no solution;
2. $\hat{w}$ has infinity solution o.w.

Situation 1 is **impossible** because both $X^\top X$ and $X^\top y$ can be represented in the column space of $X^\top$. Therefore, the optimization problem must have a solution, which may be either unique or infinite.

As an infinite set of solutions makes it difficult to determine which estimate of $\hat{w}$ to choose, we apply L_2 **regularization** to linear regression, which is commonly referred to as **Ridge Regression**. That is:

$$\mathcal{L}_{L_2} := \mathcal{L}(\hat{w}) + \boxed{\lambda \|\hat{w}\|_2^2}, \tag{1.13}$$

where $\lambda > 0$ is a hyperparameter. Notice that $\|\hat{w}\|_2^2 = \sum_{i=1}^{d+1} \hat{w}_i^2$, L_2 regularization prevents any single dimension from being assigned an excessively large weight, and encourages the model to make use of more dimensions during training.
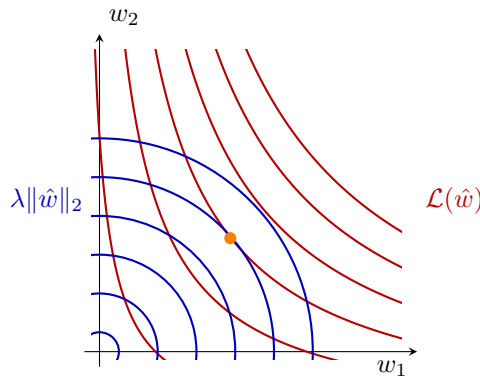

Figure 1.1: Illustration of L_2 regularization. The contours represent level sets of the regularized loss $\mathcal{L}(\hat{w}) + \lambda \|\hat{w}\|_2^2$, which take the form of concentric ellipses (circle in the plot).

During ridge regression, we minimize the $\mathcal{L}_{L_2}$:

$$\underset{\hat{w}}{\text{argmin}} \, (y - X\hat{w})^\top (y - X\hat{w}) + \lambda \hat{w}^\top \hat{w} \tag{1.14}$$

The optimization stops when:

$$\frac{\partial \mathcal{L}_{L_2}}{\partial \hat{w}} = -2X^\top y + 2X^\top X\hat{w} + 2\lambda I\hat{w} = 0 \tag{1.15}$$

$$\Rightarrow (X^\top X + \lambda I)\hat{w} = X^\top y \tag{1.16}$$

> **Proposition 1.2**
>
> $X^\top X + \lambda I$ *always invertible.* ♡

Proof Since $X^\top X$ is a real symmetric matrix, we have the eigen-decomposition $X^\top X = U \Lambda U^\top$, where $\Lambda = \text{diag}(\lambda_1, \cdots, \lambda_{d+1})$. Moreover, as $X^\top X \succeq 0$ (positive semi-definite), it follows that $\forall i \in [d+1]$, $\lambda_i \geq 0$. Note that:

$$\lambda I = \lambda U U^\top \tag{1.17}$$

since U is an orthogonal matrix. Hence:

$$X^\top X + \lambda I = U(\Lambda + \lambda I)U^\top \tag{1.18}$$

For all $i \in [d + 1]$, we have:

$$\lambda_i + \lambda > \lambda_i \geq 0 \tag{1.19}$$

Thus, $X^\top X + \lambda I$ is a full-rank matrix.

Remark Numerical issues may still occur even if $X^\top X$ is full rank (e.g., when eigenvalues λ_k are close to zero). The L_2 regularization factor λ mitigates this issue by shifting the eigenvalues upward, thereby improving numerical stability during training.

Another regularization method often used is L_1 regularization, where the loss function is defined as:

$$\mathcal{L}_{L_1} := \mathcal{L}(\hat{w}) + \boxed{\lambda \|\hat{w}\|_1} \tag{1.20}$$

L_1 regularization can induce sparsity in $\hat{w}$, which works in contrast to L_2 regularization. Specifically, L_1 regularization encourages the model to rely on only a small subset of input features, effectively performing **feature selection**.

Linear regression with L_1 regularization is called **Lasso Regression** (Least Absolute Shrinkage and Selection Operator).
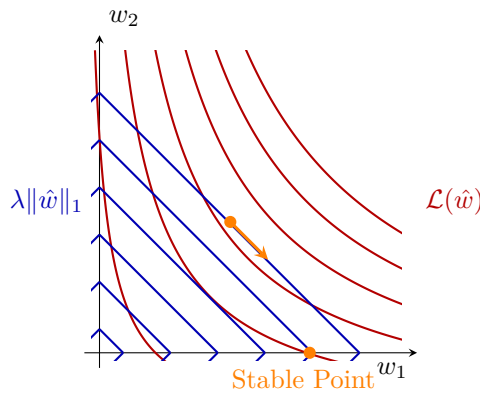

Figure 1.2: Illustration of L_1 regularization. The contours represent level sets of the regularized loss $\mathcal{L}(\hat{w}) + \lambda \|\hat{w}\|_1$, which take the form of nested diamonds (squares rotated by 45° in the plot).

1.3 Geomeric View of LR

Ideally, we would like to solve $X\hat{w} = y$. If y lies on the hypersurface

$$\mathcal{M}(X) := \mathrm{Span}(X) = \{Xw : w \in \mathbb{R}^{d+1}\} \subset \mathbb{R}^n \tag{1.21}$$

then the equation admits an exact solution. In most cases, however, $y \notin \mathcal{M}(X)$, so no exact solution exists. Nevertheless, we can always find an estimator $\hat{w}$ such that $\mathcal{P}_{\mathcal{M}(X)} y = X\hat{w}$, where $\mathcal{P}_{\mathcal{M}(X)}$ denotes the orthogonal projection onto the hypersurface $\mathcal{M}(X)$.

> **Proposition 1.3**
>
> $$\hat{y} = X\hat{w} \quad \Longrightarrow \quad \hat{w} \text{ is solution to LR.} \tag{1.22}$$

Proof

$$
\begin{aligned}
y - \hat{y} \perp \mathcal{M}(X) &\quad \Longrightarrow \quad y - X\hat{w} \perp \mathcal{M}(X) \\
&\quad \Longrightarrow \quad X^\top(y - X\hat{w}) = 0 \quad \Longrightarrow \quad \hat{w} = (X^\top X)^{-1} X^\top y
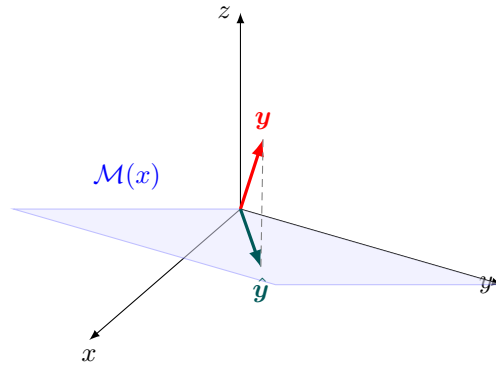\end{aligned}
$$

Figure 1.3: Orthogonal projection interpretation of linear regression. The predicted vector $X\hat{w}$ is obtained as the projection of y onto the hypersurface $\mathcal{M}(X) = \{Xw : w \in \mathbb{R}^{d+1}\}$, which is a linear subspace in the classical case.