CSC413: Homework 1

Tianyu Du (1003801647)

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1 Hard-Coding Networks

1.1 Verify Sort

Soln. The first layer performs pairwise comparison to construct indicators $\mathbb{1}\{x_1 \leq x_2\}$, $\mathbb{1}\{x_2 \leq x_3\}$, and $\mathbb{1}\{x_3 \leq x_4\}$. The second layer performs an all() operation on indicators from the previous layer.

$$\mathbf{W}^{(1)} = \begin{pmatrix} -1 & 1 & 0 & 0 \\ 0 & -1 & 1 & 0 \\ 0 & 0 & -1 & 1 \end{pmatrix} \tag{1.1}$$

$$\mathbf{b}^{(1)} = \begin{pmatrix} 0 & 0 & 0 \end{pmatrix} \tag{1.2}$$

So that

$$\varphi(\mathbf{h}) = \varphi(\mathbf{W}^{(1)}\mathbf{x} + \mathbf{b}^{(1)}) = \varphi\begin{pmatrix} x_2 - x_1 \\ x_3 - x_2 \\ x_4 - x_3 \end{pmatrix} = \begin{pmatrix} \mathbb{1}\{x_2 \ge x_1\} \\ \mathbb{1}\{x_3 \ge x_2\} \\ \mathbb{1}\{x_4 \ge x_3\} \end{pmatrix}$$
(1.3)

$$\mathbf{w}^{(2)} = \begin{pmatrix} 1 & 1 & 1 \end{pmatrix} \tag{1.4}$$

$$b^{(2)} = -0.5 (1.5)$$

Such that y = 1 if and only if all components of **h** are ones, i.e., the list is sorted.

1.2 Perform Sort

Proof. For this section, I am implementing a feedforward neural network performing bubble sort. Note that given $\mathbf{x} = (x_1, x_2, x_3, x_4)$, the **swapping** operation can be conducted by multiplying \mathbf{x} by a matrix. For instance, the following $\mathbf{W}^{(12)}$ swaps the first and second elements.

$$\underbrace{\begin{pmatrix} 0 & 1 & 0 & 0 \\ 1 & 0 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & 1 \end{pmatrix}}_{\mathbf{W}^{(12)}} (x_1, x_2, x_3, x_4)^T = (x_2, x_1, x_3, x_4)^T \tag{1.6}$$

Let I_4 denote the identity matrix. Define the following neural network:

$$\mathbf{h}_1 := \mathbb{1}\{x_1 \le x_2\}\mathbf{I}_4\mathbf{x} + \mathbb{1}\{x_1 > x_2\}\mathbf{W}^{(12)}\mathbf{x}$$
(1.7)

$$\mathbf{h}_2 = \mathbb{1}\{x_2 \le x_3\}\mathbf{I}_4\mathbf{h}_1 + \mathbb{1}\{x_2 > x_3\}\mathbf{W}^{(23)}\mathbf{h}_1$$
(1.8)

$$\mathbf{h}_3 = \mathbb{1}\{x_3 \le x_4\}\mathbf{I}_4\mathbf{h}_2 + \mathbb{1}\{x_3 > x_4\}\mathbf{W}^{(34)}\mathbf{h}_2$$
(1.9)

$$\mathbf{h}_4 = \mathbb{1}\{x_1 \le x_2\}\mathbf{I}_4\mathbf{h}_3 + \mathbb{1}\{x_1 > x_2\}\mathbf{W}^{(12)}\mathbf{h}_3$$
(1.10)

$$\mathbf{h}_5 = \mathbb{1}\{x_2 \le x_3\}\mathbf{I}_4\mathbf{h}_4 + \mathbb{1}\{x_2 > x_3\}\mathbf{W}^{(23)}\mathbf{h}_4$$
(1.11)

$$\mathbf{y}(\mathbf{x}) = \mathbb{1}\{x_1 \le x_2\}\mathbf{I}_4\mathbf{h}_5 + \mathbb{1}\{x_1 > x_2\}\mathbf{W}^{(12)}\mathbf{h}_5$$
(1.12)

The above neural network can also be implemented using a recurrent structure, in which hidden states record the number of pairwise comparisons made.9

1.3 Universal Approximation Theorem

1.3.1

Soln. To avoid over-using of notations, let $\varphi(y) := \mathbb{1}\{y > 0\}$ denote the activation function.

$$n = 2 \tag{1.13}$$

$$\mathbf{W}_0 = (1, -1) \tag{1.14}$$

$$\mathbf{b}_0 = (-a, b) \tag{1.15}$$

$$\mathbf{W}_1 = (1, 1) \tag{1.16}$$

$$\mathbf{b}_1 = -0.5 \tag{1.17}$$

Justification:

$$\varphi(\mathbf{h}) = \varphi((x - a, b - x)) \tag{1.18}$$

$$= (1\{x - a > 0\}, 1\{b - x > 0\})$$
(1.19)

$$= (\mathbb{1}\{x > a\}, \mathbb{1}\{x < b\}) \tag{1.20}$$

$$\varphi(\mathbf{W}_1 \varphi(\mathbf{h}) + \mathbf{b}_1) = \mathbb{1} \{ \mathbb{1} \{ x > a \} + \mathbb{1} \{ x < b \} - 0.5 \}$$
(1.21)

$$= 1\{x > a\} \land 1\{x < b\} \tag{1.22}$$

$$= 1{a < x < b} \tag{1.23}$$

Soln. Let $\delta \in (0,1)$ denote the ratio parameter, a higher value of δ results in a finer approximation is. In this example, take $\delta = \frac{9}{10}$.

Without loss of generality, assume the region I on which function f is defined on to be symmetric across zero.

Let I = [-1, 1], given f is symmetric, $f(-\delta) = f(\delta)$.

Define:

$$\hat{f}_1(x) = \hat{f}_0(x) + g(f(\delta), -\delta, \delta, x)$$

$$(1.24)$$

Note that

$$||f - \hat{f}_1|| = \int_{-1}^{1} |f(x) - \hat{f}_1(x)| dx$$
 (1.25)

$$= \int_{-1}^{-\delta} |f(x)| dx + \int_{-\delta}^{\delta} |f(x) - \hat{f}_1(x)| dx + \int_{\delta}^{1} |f(x)| dx$$
 (1.26)

Given that $\forall x \in (-\delta, \delta), \ f(x) > f(-\delta) = f(\delta) > 0$, it follows

$$\int_{-\delta}^{\delta} |f(x) - \hat{f}_1(x)| dx = \int_{-\delta}^{\delta} f(x) - \hat{f}_1(x) dx$$
 (1.27)

$$= \int_{-\delta}^{\delta} f(x) dx - \int_{-\delta}^{\delta} \hat{f}_1(x) dx$$
 (1.28)

Also, $\int_{-\delta}^{\delta} \hat{f}_1(x) \ dx > 0$ provided $\delta \neq 0$. Therefore,

$$\int_{-\delta}^{\delta} \left| f(x) - \hat{f}_1(x) \right| dx < \int_{-\delta}^{\delta} f(x) dx \tag{1.29}$$

$$= \int_{-\delta}^{\delta} |f(x)| \ dx \tag{1.30}$$

Therefore,

$$||f - \hat{f}_1|| = \int_{-1}^{-\delta} |f(x)| dx + \int_{-\delta}^{\delta} |f(x) - \hat{f}_1(x)| dx + \int_{\delta}^{1} |f(x)| dx$$
 (1.31)

$$<\int_{-1}^{-\delta} |f(x)| dx + \int_{-\delta}^{\delta} |f(x)| dx + \int_{\delta}^{1} |f(x)| dx$$
 (1.32)

$$= \int_{-1}^{1} |f(x) - 0| \ dx \tag{1.33}$$

$$= \int_{-1}^{1} \left| f(x) - \hat{f}_0(x) \right| dx \tag{1.34}$$

$$= ||f(x) - \hat{f}_0(x)|| \tag{1.35}$$

Therefore,

$$||f(x) - \hat{f}_1(x)|| < ||f(x) - \hat{f}_0(x)||$$
 (1.36)

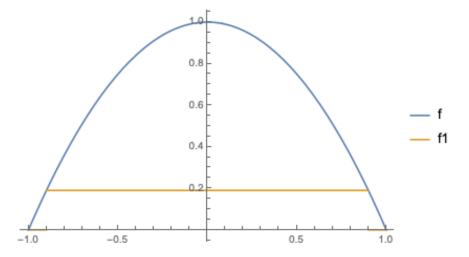


Figure 1.1: Approximated Result

Soln. Algorithm:

(i) Divide I = [-1, 1] into N + 2 sub-intervals with equal length, such that

$$I_i := \left[-1 + \frac{i}{N+2}, -1 + \frac{i+1}{N+2} \right] \quad \forall \ i \in \{1, 2, \dots, N\}$$
 (1.37)

Note that the first and last sub-intervals are not used to construct g_i .

(ii) For each i, define

$$h_i := \min_{x \in I_i} f(x) \tag{1.38}$$

$$a_i := -1 + \frac{i}{N+2} \tag{1.39}$$

$$b_i := -1 + \frac{i+1}{N+2} \tag{1.40}$$

Because $f(x) \ge 0 \ \forall x \in I$.

By the definition of $g_i(x)$, it can be shown that¹

$$f(x) \ge f_i(x) \ \forall i \in \{1, 2, \dots, N\} \ \forall x \in \bigcup_{i=1}^N (a_i, b_i)$$
 (1.41)

Further,

$$f(x) = f_i(x) \quad \forall i \in \{1, 2, \dots, N\} \ \forall x \in \left[-1, -1 + \frac{1}{N+2}\right] \bigcup \left(1 - \frac{1}{N+2}, 1\right]$$
 (1.42)

 $^{^{1}}$ I am excluding those boundary points between consecutive sub-intervals, because at those points, the value of f_{i} spikes due to duplicate counts of indicator functions. However, while doing integral, this does not matter as the set of boundary points has measure zero.

Define

$$\mathcal{K} := \left[-1, -1 + \frac{1}{N+2} \right) \bigcup \left(1 - \frac{1}{N+2}, 1 \right] \bigcup \left(\bigcup_{i=1}^{N} (a_i, b_i) \right)$$
 (1.43)

Note that the set $I \setminus \mathcal{K}$ consists of all boundary points between consecutive sub-intervals. There are only finitely many such points, therefore $I \setminus \mathcal{K}$ has measure zero, and

$$\int_{I} \left| f(x) - \hat{f}_{i}(x) \right| dx = \int_{\mathcal{K}} \left| f(x) - \hat{f}_{i}(x) \right| dx \tag{1.44}$$

And I've shown that for every i and every $x \in \mathcal{K}$, $f(x) \geq f_i(x)$. Consequently,

$$\int_{I} \left| f(x) - \hat{f}_{i}(x) \right| dx = \int_{\mathcal{K}} \left| f(x) - \hat{f}_{i}(x) \right| dx \text{ (removing measure zero set.)}$$
(1.45)

$$= \int_{\mathcal{K}} f(x) - \hat{f}_i(x) \, dx \tag{1.46}$$

$$= \int_{I} f(x) - \hat{f}_{i}(x) dx \text{ (adding back the measure zero set.)} \quad (\dagger)$$
 (1.47)

Define $\hat{f}_0(x) = 0$ and let $i \in \{1, 2, \dots, N\}$,

$$||f - \hat{f}_{i+1}|| = \int_{-1}^{1} |f(x) - \hat{f}_{i+1}(x)| dx$$
(1.48)

$$= \int_{-1}^{1} f(x) - \hat{f}_{i+1}(x) \, dx \text{ by (†)}$$
(1.49)

$$= \int_{-1}^{-1\frac{i+1}{N}} f(x) - \hat{f}_{i+1}(x) dx + \int_{-1+\frac{i+1}{N}}^{-1+\frac{i+2}{N}} f(x) - \hat{f}_{i+1}(x) dx + \int_{-1+\frac{i+2}{N}}^{1} f(x) - \hat{f}_{i+1}(x) dx$$

$$(1.50)$$

Further, by construction, $\hat{f}_{i+1}(x) = \hat{f}_i(x) \ \forall x \notin [a_{i+1}, b_{i+1}]$. Therefore,

$$||f - \hat{f}_{i+1}|| = \int_{-1}^{a_{i+1}} f(x) - \hat{f}_i(x) \, dx + \int_{a_{i+1}}^{b_{i+1}} f(x) - \hat{f}_{i+1}(x) \, dx + \int_{b_{i+1}}^{1} f(x) - \hat{f}_i(x) \, dx$$

$$(1.51)$$

$$= \int_{-1}^{a_{i+1}} f(x) - \hat{f}_i(x) dx + \int_{a_{i+1}}^{b_{i+1}} f(x) - \hat{f}_i(x) - g(h_{i+1}, a_{i+1}, b_{i+1}, x) dx + \int_{b_i}^{1} f(x) - \hat{f}_i(x) dx$$

$$(1.52)$$

$$= \int_{-1}^{a_{i+1}} f(x) - \hat{f}_i(x) dx + \int_{a_{i+1}}^{b_{i+1}} f(x) - \hat{f}_i(x) dx + \int_{b_{i+1}}^{1} f(x) - \hat{f}_i(x) dx - \int_{a_{i+1}}^{b_{i+1}} g(h_{i+1}, a_{i+1}, b_{i+1}, x) dx$$

$$(1.53)$$

$$= \int_{-1}^{1} f(x) - \hat{f}_i(x) dx - \int_{a_{i+1}}^{b_{i+1}} g(h_{i+1}, a_{i+1}, b_{i+1}, x) dx$$
(1.54)

$$= ||f - \hat{f}_i|| - \int_{a_{i+1}}^{b_{i+1}} g(h_{i+1}, a_{i+1}, b_{i+1}, x) dx$$
(1.55)

Note that for every i, for every $x \in [a_i, b_i]$, $g(h_i, a_i, b_i, x) > 0$. Therefore, $\int_{a_i}^{b_i} g(h_i, a_i, b_i, x) dx > 0$. Hence,

$$||f - \hat{f}_{i+1}|| = ||f - \hat{f}_{i}|| - \int_{a_{i+1}}^{b_{i+1}} g(h_{i+1}, a_{i+1}, b_{i+1}, x) dx$$

$$> ||f - \hat{f}_{i}||$$

$$(1.56)$$

$$> ||f - \hat{f}_i|| \tag{1.57}$$

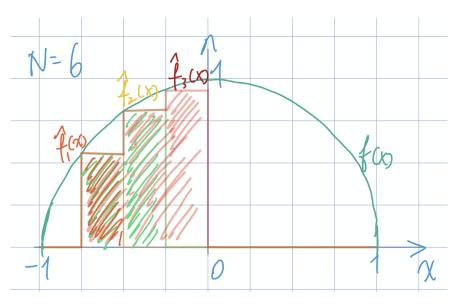


Figure 1.2: Approximated Results for ${\cal N}=6$

1.3.4

Soln. Not required.

2 Backprop

2.1 Computational Graph

2.1.1

Soln.

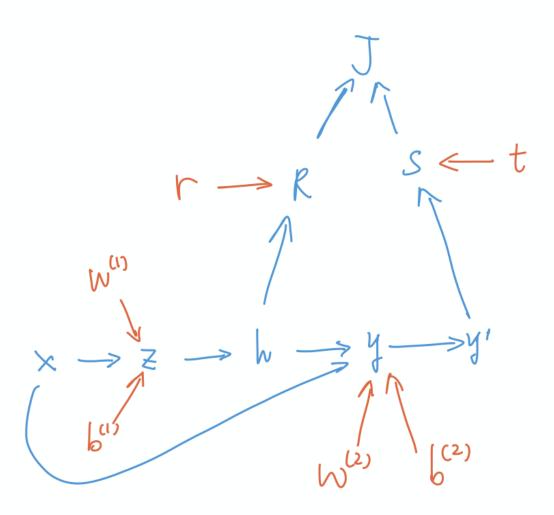


Figure 2.1: Computational Graph

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2.1.2

Soln.

$$\overline{\mathbf{x}} = \overline{\mathbf{z}} \frac{\partial \mathbf{z}}{\partial \mathbf{x}} \tag{2.1}$$

$$= \overline{\mathbf{z}}\mathbf{W}^{(1)} \tag{2.2}$$

$$= \overline{\mathbf{h}} \frac{\partial \mathbf{h}}{\partial \mathbf{z}} \mathbf{W}^{(1)} \tag{2.3}$$

$$= \overline{\mathbf{h}} \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)} \tag{2.4}$$

$$= \left(\overline{\mathcal{R}}\frac{\partial \mathcal{R}}{\partial \mathbf{h}} + \overline{\mathbf{y}}\frac{\partial \mathbf{y}}{\partial \mathbf{h}}\right) \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)}$$
(2.5)

$$= \left(\overline{\mathcal{R}}\mathbf{r}^T + \overline{\mathbf{y}}\mathbf{W}^{(2)}\right) \mathbb{1}\{\mathbf{z} \ge 0\}\mathbf{W}^{(1)}$$
(2.6)

$$= \left(\mathbf{r}^T + \overline{\mathbf{y}'} \frac{\partial \mathbf{y}'}{\partial \mathbf{y}} \mathbf{W}^{(2)}\right) \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)}$$
(2.7)

$$= \left(\mathbf{r}^T + \overline{\mathbf{y}'} \mathtt{softmax}'(\mathbf{y}) \mathbf{W}^{(2)}\right) \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)}$$
(2.8)

$$= \left(\mathbf{r}^T + \overline{\mathcal{S}} \frac{\partial \mathcal{S}}{\partial \mathbf{y}'} \operatorname{softmax}'(\mathbf{y}) \mathbf{W}^{(2)}\right) \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)}$$
(2.9)

$$= \left(\mathbf{r}^T + \mathbf{e}_k \operatorname{softmax}'(\mathbf{y}) \mathbf{W}^{(2)}\right) \mathbb{1}\{\mathbf{z} \ge 0\} \mathbf{W}^{(1)}$$
(2.10)

where \mathbf{e}_k denotes the one-hot vector in \mathbb{R}^M in which the k^{th} element is one.

2.2 Vector-Jacobean Product (VJPs)

2.2.1

Soln.

$$\mathbf{f}(\mathbf{x}) := \mathbf{v}\mathbf{v}^T\mathbf{x} \tag{2.11}$$

$$\implies J = \frac{d}{d\mathbf{x}}\mathbf{f}(\mathbf{x}) = \mathbf{v}\mathbf{v}^T \tag{2.12}$$

$$\implies J = \begin{pmatrix} 1 & 2 & 3 \\ 2 & 4 & 6 \\ 3 & 6 & 9 \end{pmatrix} \tag{2.13}$$

2.2.2

Soln. Time cost: n^2 .

Memory cost: $\frac{(n+1)n}{2}$ (only need to store elements above the diagonal (including the diagonal), because J is symmetric).

2.2.3

Soln.

$$J^T \mathbf{y} = \mathbf{v} \mathbf{v}^T \mathbf{y} \tag{2.14}$$

$$= \mathbf{v}(\mathbf{v}^T \mathbf{y}) \tag{2.15}$$

$$=\underbrace{\mathbf{v}}_{\text{step 2}}^{\text{step 1}} \quad (\dagger) \tag{2.16}$$

where the inner product operation has time cost n and its output takes memory cost 1. Let $\alpha = \langle \mathbf{v}, \mathbf{y} \rangle \in \mathbb{R}$, the vector scaler multiplication operation \mathbf{v}_{α} has time cost n and its output has memory cost n.

The overall time cost for (\dagger) is therefore 2n and memory cost is n+1.

Linear Regression $\mathbf{3}$

3.1Driving the Gradient

Soln.

$$\frac{d}{d\hat{\mathbf{w}}} \frac{1}{n} (X\hat{\mathbf{w}} - \mathbf{t})^2 = \frac{d}{d\hat{\mathbf{w}}} \frac{1}{n} ||X\hat{\mathbf{w}} - \mathbf{t}||_2^2$$
(3.1)

$$= \frac{2}{n} (X\hat{\mathbf{w}} - \mathbf{t})^T X \tag{3.2}$$

$$= \frac{2}{n} (X\hat{\mathbf{w}} - \mathbf{t})^T X$$

$$\implies \nabla_{\mathbf{w}} \mathcal{J} = \frac{2}{n} X^T (X\hat{\mathbf{w}} - \mathbf{t})$$
(3.2)

3.2 **Under-parameterized Model**

3.2.1

Soln. Assume d < n so that X^TX is invertible. The gradient descent algorithm converges when the gradient equals zero:

$$\frac{2}{n}(X\hat{\mathbf{w}} - \mathbf{t})^T X = 0 \tag{3.4}$$

$$\implies (X\hat{\mathbf{w}} - \mathbf{t})^T X = 0 \tag{3.5}$$

$$\implies X^T(X\hat{\mathbf{w}} - \mathbf{t}) = 0^T \tag{3.6}$$

$$\implies X^T X \hat{\mathbf{w}} - X^T \mathbf{t} = 0^T \tag{3.7}$$

$$\implies X^T X \hat{\mathbf{w}} = X^T \mathbf{t} \tag{3.8}$$

$$\implies \hat{\mathbf{w}} = (X^T X)^{-1} X^T \mathbf{t} \tag{3.9}$$

3.2.2

Soln. Let $\mathbf{x} \in \mathbb{R}^d$, note that $(X^TX)^{-1}$ is symmetric. Assuming target \mathbf{t} is generated by a linear process, then $\mathbf{t} = X\mathbf{w}^*$. Immediately, $\mathbf{t}^T = \mathbf{w}^{*T}X^T$.

$$(\mathbf{w}^{*T}\mathbf{x} - \hat{\mathbf{w}}^T\mathbf{x})^2 = (\mathbf{w}^{*T}\mathbf{x} - [(X^TX)^{-1}X^T\mathbf{t}]^T\mathbf{x})^2$$
(3.10)

$$= (\mathbf{w}^{*T}\mathbf{x} - \mathbf{t}^T X (X^T X)^{-1} \mathbf{x})^2$$
(3.11)

$$= (\mathbf{w}^{*T}\mathbf{x} - \mathbf{w}^{*T}X^TX(X^TX)^{-1}\mathbf{x})^2$$
(3.12)

$$= (\mathbf{w}^{*T}\mathbf{x} - \mathbf{w}^{*T}\mathbf{x})^2 \tag{3.13}$$

$$=0 (3.14)$$

3.3 Over-parameterized Model: 2D Example

3.3.1

Soln. To minimize the empirical risk minimizer,

$$\min_{w_1, w_2} (w_1 x_1 + w_2 x_2 - t_1)^2 \tag{3.15}$$

$$\min_{w_1, w_2} (w_1 x_1 + w_2 x_2 - t_1)^2$$
equivalently,
$$\min_{w_1, w_2} (2w_1 + w_2 - 2)^2$$
(3.15)

Any pair of (w_1, w_2) satisfying

$$2w_1 + w_2 - 2 = 0 \quad (\dagger) \tag{3.17}$$

attains the minimum level of empirical risk (zero). Equivalently, any $\hat{\mathbf{w}}$ on the line

$$\hat{\mathbf{w}} = \begin{pmatrix} 0 \\ 2 \end{pmatrix} + t \begin{pmatrix} 1 \\ -2 \end{pmatrix} \text{ for } t \in \mathbb{R}$$
 (3.18)

satisfies (†). Therefore, there are infinitely many empirical risk minimizers. Equivalently, the collection of solution is

$$w_2 = -2w_1 + 2 \tag{3.19}$$

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Soln. From the fist part of this question we know that

$$\nabla_{\mathbf{w}} \mathcal{J} = \frac{2}{n} (X \hat{\mathbf{w}} - \mathbf{t})^T X \tag{3.20}$$

$$= \frac{2}{1} \left[\begin{pmatrix} 2 \\ 1 \end{pmatrix} \cdot \begin{pmatrix} 0 \\ 0 \end{pmatrix} - 2 \right] \begin{pmatrix} 2 \\ 1 \end{pmatrix} \tag{3.21}$$

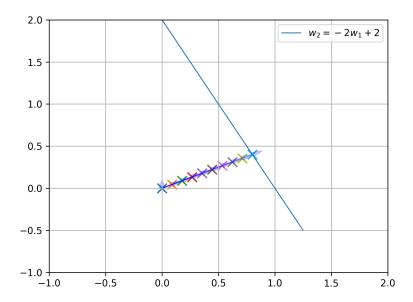
$$= \begin{pmatrix} -4\\ -2 \end{pmatrix} \tag{3.22}$$

The unit-norm gradient is

$$\widehat{\nabla_{\mathbf{w}}} \mathcal{J} = \begin{pmatrix} -\frac{2\sqrt{5}}{5} \\ -\frac{\sqrt{5}}{5} \end{pmatrix} \tag{3.23}$$

The direction (gradient) does not change along the trajectory. Ultimately, the gradient descend algorithm converges to

$$\hat{\mathbf{w}}^* = \left(\frac{4}{5}, \frac{2}{5}\right) \tag{3.24}$$



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Soln. Let $\hat{\mathbf{w}}^*$ denote the solution found using gradient descent. Note that the line of solution can be written parametrically as

$$\hat{\mathbf{w}}(t) = \begin{pmatrix} 0 \\ 2 \end{pmatrix} + t \begin{pmatrix} 1 \\ -2 \end{pmatrix} \tag{3.25}$$

The path of gradient descent starts from S = 0 and the path is perpendicular to the line of solutions $\{\hat{\mathbf{w}}(t):t\in\mathbb{R}\}$. By Pythagorean theorem, we know that the shortest path from a fixed point x to a line ℓ is the perpendicular line x and some point y on ℓ . Meanwhile, y is the point on ℓ nearest to x. The path of gradient descent is exactly such a shortest path. Therefore, $\hat{\mathbf{w}}^*$ is the solution nearest to $\mathbf{0}$. In another word, $\hat{\mathbf{w}}$ is solution with smallest Euclidean among all solutions in \mathcal{S} .

3.4 Overparameterized Model: General Case

3.4.1

Proof. Note that the solution reached by gradient descent \mathbf{w}^* can be written as a linear combination of rows of X:

$$X = \begin{pmatrix} x_{11} & \cdots & x_{1d} \\ \vdots & \ddots & \vdots \\ x_{n1} & \cdots & x_{nd} \end{pmatrix}$$
 (3.26)

$$\mathbf{w}^* = r_1 \begin{pmatrix} x_{11} \\ \vdots \\ x_{1d} \end{pmatrix} + \dots + r_n \begin{pmatrix} x_{n1} \\ \vdots \\ x_{nd} \end{pmatrix}$$
(3.27)

$$= \begin{pmatrix} x_{11} & \cdots & x_{1d} \\ \vdots & \ddots & \vdots \\ x_{n1} & \cdots & x_{nd} \end{pmatrix}^T \begin{pmatrix} r_1 \\ \vdots \\ r_n \end{pmatrix}$$
(3.28)

$$=X^{T}\mathbf{r}\tag{3.29}$$

The original minimization problem in \mathbb{R}^d (choosing the optimal $\mathbf{w} \in \mathbb{R}^d$) can be reduced to a minimization problem in \mathbb{R}^n (choosing the optimal $\mathbf{r} \in \mathbb{R}^n$).

The gradient descent algorithm converges if and only if zero gradient is encountered, which is characterized by the following condition:

$$\frac{d}{d\mathbf{r}}\frac{1}{n}||XX^T\mathbf{r} - \mathbf{t}||_2^2 = 0 \tag{3.30}$$

$$\frac{d}{d\mathbf{r}} \frac{1}{n} ||XX^T \mathbf{r} - \mathbf{t}||_2^2 = 0$$

$$\Rightarrow \frac{2}{n} (XX^T \mathbf{r} - \mathbf{t})^T XX^T = 0$$
(3.30)

$$\implies XX^T(XX^T\mathbf{r} - \mathbf{t}) = 0$$
 (take transpose on both sides) (3.32)

$$\implies XX^TXX^T\mathbf{r} = XX^T\mathbf{t} \tag{3.33}$$

Because d > n, so $rank(XX^T) = n$ and it is invertible. Therefore,

$$(XX^{T})^{-1}XX^{T}XX^{T}\mathbf{r} = (XX^{T})^{-1}XX^{T}\mathbf{t}$$
(3.34)

$$\implies XX^T\mathbf{r} = \mathbf{t} \tag{3.35}$$

$$\implies \mathbf{r}^* = (XX^T)^{-1}\mathbf{t} \tag{3.36}$$

$$\implies \mathbf{w}^* = X^T (XX^T)^{-1} \mathbf{t} \tag{3.37}$$

where the solution is uniquely determined.

3.4.2

Proof. Let $\hat{\mathbf{w}}_1$ be another zero loss weight, then

$$X\hat{\mathbf{w}}_1 = \mathbf{t} \tag{3.38}$$

Immediately,

$$\hat{\mathbf{w}}_1^T X^T = \mathbf{t}^T \tag{3.39}$$

Evaluating

$$(\hat{\mathbf{w}} - \hat{\mathbf{w}}_1)^T \hat{\mathbf{w}} = ||\hat{\mathbf{w}}||_2^2 - \hat{\mathbf{w}}_1^T \hat{\mathbf{w}}$$

$$(3.40)$$

$$= (X^T (XX^T)^{-1} \mathbf{t})^T (X^T (XX^T)^{-1} \mathbf{t}) - \hat{\mathbf{w}}_1^T \hat{\mathbf{w}}$$
(3.41)

$$= \mathbf{t}^T (XX^T)^{-1} X (X^T (XX^T)^{-1} \mathbf{t}) - \hat{\mathbf{w}}_1^T \hat{\mathbf{w}}$$
(3.42)

$$= \mathbf{t}^{T} (XX^{T})^{-1} \mathbf{t} - \hat{\mathbf{w}}_{1}^{T} X^{T} (XX^{T})^{-1} \mathbf{t}$$
(3.43)

$$= \mathbf{t}^T (XX^T)^{-1} \mathbf{t} - \mathbf{t}^T (XX^T)^{-1} \mathbf{t}$$
(3.44)

$$= 0 \quad (\dagger) \tag{3.45}$$

Because $\hat{\mathbf{w}}_1$ is chosen arbitrarily, let \mathcal{S} denote the space of zero-loss weights. (†) suggests the optimal solution found by gradient descent $\hat{\mathbf{w}} \perp \mathcal{S}$.

Let $\hat{\mathbf{w}}$ denote the solution found by gradient descent.

Let $\hat{\mathbf{w}}_{\alpha} \in \mathcal{S}$ and $\mathbf{v} = \hat{\mathbf{w}}_{\alpha} - \hat{\mathbf{w}}$.

$$||\hat{\mathbf{w}}_{\alpha}||_2^2 = ||\hat{\mathbf{w}} + \mathbf{v}||_2^2 \tag{3.46}$$

$$= ||\hat{\mathbf{w}}||_{2}^{2} + 2\mathbf{v} \cdot \hat{\mathbf{w}} + ||\mathbf{v}||_{2}^{2} \tag{3.47}$$

$$= ||\hat{\mathbf{w}}||_{2}^{2} + 2(\hat{\mathbf{w}}_{\alpha} - \hat{\mathbf{w}}) \cdot \hat{\mathbf{w}} + ||\mathbf{v}||_{2}^{2}$$
(3.48)

$$= ||\hat{\mathbf{w}}||_2^2 + ||\mathbf{v}||_2^2 \text{ by (†)}$$
(3.49)

$$\geq ||\hat{\mathbf{w}}||_2^2 \tag{3.50}$$

Therefore,

$$||\hat{\mathbf{w}}_{\alpha}||_{2}^{2} \ge ||\hat{\mathbf{w}}||_{2}^{2} \quad \forall \ \hat{\mathbf{w}}_{\alpha} \in \mathcal{S} \tag{3.51}$$

And $\hat{\mathbf{w}}$ is the element with smallest Euclidean norm among all zero-loss solutions in \mathcal{S} .

3.5 Benefit of Overparameterization

3.5.1

Soln. Fitting function implementation:

```
def fit_poly(X, d,):
    X_expand = poly_expand(X, d=d, poly_type = poly_type)
    if d > n: # Over-parameterized case
    W = X_expand.T @ np.linalg.inv(X_expand @ X_expand.T) @ t
    else: # Under parameterized case.
    W = np.linalg.inv(X_expand.T @ X_expand) @ X_expand.T @ t
    return W
```

Lines fitted and the corresponding losses using both legendre and chebyshev polynomials are presented in figures below. In both experiments, over-parameterization does **not** always lead to overfitting. In fact, models with a super high degree (say 100) actually outperformed models with low degrees.

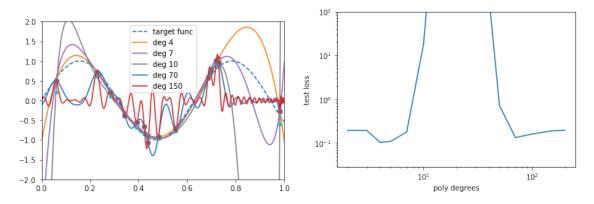


Figure 3.1: chebyshev results

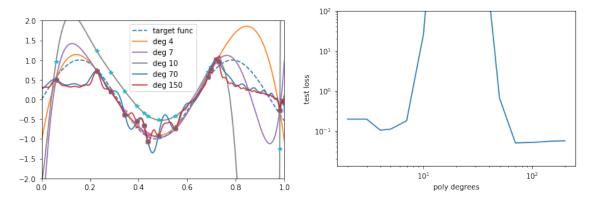


Figure 3.2: legendre results

3.5.2

Soln. Not required.