Training One Neural Layer

Deep Learning (ENGG*6600*01)

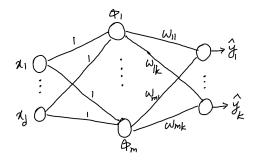
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Introduction

- In the lecture of "Training one neuron", we were introduced to the neural networks with only one layer and only one neuron.
- In this lecture, we are introduced to the neural networks with only one [learnable] layer but possibly multiple neurons.
- Two of these networks are Radial Basis Function (RBF) network and Self-Organizing Map (SOM).

- A Radial Basis Function (RBF) network was first proposed in 1988 [1, 2].
- It has two layers but the first layer has fixed weights equal to one. The second layer has learnable weights.
- The first layer connects the data $\mathbf{x} \in \mathbb{R}^d$ to m basis kernel functions $\{\phi_i(\mathbf{x})\}_{i=1}^m$.
- The second layer connects the basis functions $\{\phi_i(\mathbf{x})\}_{i=1}^m$ to the output neurons $\{\widehat{y}_j\}_{j=1}^k$.
- The weight w_{ij} denotes the weight connecting $\phi_i(\mathbf{x})$ to \hat{y}_i .



The basis functions can be various kernel functions such as:

Gaussian distribution:
$$\phi_i(\mathbf{x}) = e^{-\frac{\|\mathbf{x}_i - \mu_i\|_2^2}{2\sigma_i^2}}$$
, (1)

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Logistic (sigmoid) function: $\phi_i(\mathbf{x}) = \frac{1}{1 + e^{-\frac{\|\mathbf{x}_i - \boldsymbol{\mu}_i\|_2^2}{2\sigma_i^2}}}$, (2)

where $\mu_i \in \mathbb{R}^d$ and $\sigma_i^2 \in \mathbb{R}$ are the mean and variance for $\phi_i(\mathbf{x})$.

• At the first step, the means $\{\mu_i\}_{i=1}^m$ are found by applying a clustering method, such as K-means, on the training data with m clusters. The variances of clusters determine the variances $\{\sigma_i\}_{i=1}^m$.

- RBF networks can be considered as an additive model. An additive model, first proposed
 in [3], maps data to a space with several basis functions and then tries to learn a weighted
 average of those bases.
- In RBF, the output is obtained as:

$$\widehat{y}_j = w_{1j} \phi_1(\mathbf{x}) + \dots + w_{mj} \phi_m(\mathbf{x}) = \sum_{i=1}^m w_{ij} \phi_i(\mathbf{x}).$$
 (3)

• Consider *n* data points together in a matrix $X \in \mathbb{R}^{k \times n}$. In matrix form:

$$\widehat{\mathbf{Y}} = \mathbf{W}^{\top} \mathbf{\Phi},\tag{4}$$

where:

$$\mathbb{R}^{k \times n} \ni \widehat{\mathbf{Y}} = \begin{bmatrix} \widehat{y}_{11} & \dots & \widehat{y}_{1n} \\ \vdots & \ddots & \vdots \\ \widehat{y}_{k1} & \dots & \widehat{y}_{kn} \end{bmatrix}, \quad \mathbb{R}^{m \times k} \ni \mathbf{W} = \begin{bmatrix} w_{11} & \dots & w_{1k} \\ \vdots & \ddots & \vdots \\ w_{m1} & \dots & w_{mk} \end{bmatrix},$$

$$\mathbb{R}^{m \times n} \ni \mathbf{\Phi} = \begin{bmatrix} \phi_1(x_1) & \dots & \phi_1(x_n) \\ \vdots & \ddots & \vdots \\ \phi_m(x_1) & \dots & \phi_m(x_n) \end{bmatrix}.$$

• Least squares error minimization between the label of data $Y \in \mathbb{R}^{k \times n}$ and the output of network $\widehat{Y} \in \mathbb{R}^{k \times n}$:

$$\underset{\boldsymbol{W}}{\text{minimize}} \quad \|\boldsymbol{Y} - \widehat{\boldsymbol{Y}}\|_F^2 = \|\boldsymbol{Y} - \boldsymbol{W}^\top \boldsymbol{\Phi}\|_F^2. \tag{5}$$

Simplification of the cost function:

$$\begin{split} &\|\boldsymbol{Y} - \boldsymbol{W}^\top \boldsymbol{\Phi}\|_F^2 = \operatorname{tr} \big((\boldsymbol{Y} - \boldsymbol{W}^\top \boldsymbol{\Phi})^\top (\boldsymbol{Y} - \boldsymbol{W}^\top \boldsymbol{\Phi}) \big) = \operatorname{tr} \big((\boldsymbol{Y}^\top - \boldsymbol{\Phi}^\top \boldsymbol{W}) (\boldsymbol{Y} - \boldsymbol{W}^\top \boldsymbol{\Phi}) \big) \\ &= \operatorname{tr} (\boldsymbol{Y}^\top \boldsymbol{Y} - \boldsymbol{Y}^\top \boldsymbol{W}^\top \boldsymbol{\Phi} - \boldsymbol{\Phi}^\top \boldsymbol{W} \boldsymbol{Y} + \boldsymbol{\Phi}^\top \boldsymbol{W} \boldsymbol{W}^\top \boldsymbol{\Phi}) \\ &\stackrel{(a)}{=} \operatorname{tr} (\boldsymbol{Y}^\top \boldsymbol{Y}) - \operatorname{tr} (\boldsymbol{W}^\top \boldsymbol{\Phi} \boldsymbol{Y}^\top) - \operatorname{tr} (\boldsymbol{W} \boldsymbol{Y} \boldsymbol{\Phi}^\top) + \operatorname{tr} (\boldsymbol{W}^\top \boldsymbol{\Phi} \boldsymbol{\Phi}^\top \boldsymbol{W}), \end{split}$$

where (a) is because of the linearity and cyclic property of the trace operator.

Solving this optimization problem:

$$\frac{\partial}{\partial \mathbf{W}} \| \mathbf{Y} - \mathbf{W}^{\top} \mathbf{\Phi} \|_{F}^{2} = -\mathbf{\Phi} \mathbf{Y}^{\top} - \mathbf{\Phi} \mathbf{Y}^{\top} + 2\mathbf{\Phi} \mathbf{\Phi}^{\top} \mathbf{W} \stackrel{\text{set}}{=} \mathbf{0} \implies \mathbf{\Phi} \mathbf{\Phi}^{\top} \mathbf{W} = \mathbf{\Phi} \mathbf{Y}^{\top}$$

$$\implies \mathbf{W} = (\mathbf{\Phi} \mathbf{\Phi}^{\top})^{-1} \mathbf{\Phi} \mathbf{Y}^{\top}. \tag{6}$$

• If $\Phi\Phi^{\top}$ is a singular matrix, the pseudo-inverse can be used:

$$\mathbf{W} = (\mathbf{\Phi}\mathbf{\Phi}^{\top})^{\dagger}\mathbf{\Phi}\mathbf{Y}^{\top},\tag{7}$$

where † denotes the pseudo-inverse of matrix.

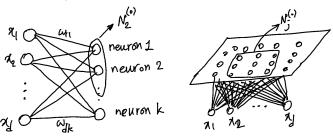
• The output of the RBF network, for either the training or test data, is:

$$\widehat{\mathbf{Y}} = \mathbf{W}^{\top} \mathbf{\Phi} = \left((\mathbf{\Phi} \mathbf{\Phi}^{\top})^{-1} \mathbf{\Phi} \mathbf{Y}^{\top} \right)^{\top} \mathbf{\Phi} = \mathbf{Y} \mathbf{\Phi}^{\top} (\mathbf{\Phi} \mathbf{\Phi}^{\top})^{-\top} \mathbf{\Phi} = \mathbf{Y} \mathbf{\Phi}^{\top} (\mathbf{\Phi} \mathbf{\Phi}^{\top})^{-1} \mathbf{\Phi}. \tag{8}$$

Self-Organizing Map

Self-Organizing Map

- Self-Organizing Map (SOM) is a neural network with one layer. It is used for unsupervised clustering, where the name "self-organizing" comes from.
- It was proposed by Teuvo Kohonen in 1982 [4]; therefore, it is also called the Kohonen network [5].
- It is one layer connecting $\mathbf{x} = [x_1, \dots, x_d]^{\top} \in \mathbb{R}^d$ to k neurons. Let w_{ij} denote the weight connecting x_i to the j-th neuron. Each neuron represents a cluster. SOM trains the weights to cluster the input data \mathbf{x} to one of the k clusters.
- The neurons can be put in 1D or 2D structure.
- Every neuron has a neighborhood around it in the 1D or 2D structure of neurons. This neighborhood is decreased gradually during the training phase. Let $\mathcal{N}_j^{(\tau)}$ denote the neighborhood of the j-th neuron at iteration τ .



Self-Organizing Map

- Step 1 of training: Initialize all weights to small random values. Set all neighborhoods $\{\mathcal{N}_j^{(0)}\}_{j=1}^k$ to half of the neuron structure grid. Set the initial learning rate $\eta^{(0)}$ to a number in range (0,1].
- Step 2 of training: Select one the input data points, $\mathbf{x} = [x_1, \dots, x_d]^\top$, and feed it to the network. Select the wining neuron \mathbf{z} by:

$$z := \arg\min_{j} \sum_{i=1}^{d} \|x_i - w_{ij}\|_2.$$
 (9)

 Step 3 of training: Update the weights of the neurons in the neighborhood of the winning neuron z:

$$w_{ij}^{(\tau+1)} := \begin{cases} w_{ij}^{(\tau)} + \eta^{(\tau)} (x_i - w_{ij}^{(\tau)}) & \text{if } j \in \mathcal{N}_z^{(\tau)} \\ w_{ij}^{(\tau)} & \text{Otherwise,} \end{cases}$$
(10)

for all $i \in \{1, ..., d\}$.

• Step 4 of training: Decrease the learning rate and the neighborhoods:

$$\eta^{(\tau+1)} := \eta^{(0)} (1 - \frac{\tau}{t}),\tag{11}$$

$$\mathcal{N}_{j}^{(\tau+1)} := \mathcal{N}_{j}^{(\tau+1)}/2, \quad \forall j \in \{1, \dots, k\},$$
 (12)

where t denotes the total number of training iterations.

• Step 5 of training: Increase τ and go to step 2.

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