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# Deep Learning Assignment 3

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## 1 General Questions

(a) Say if the first module is:

$$\max(W_1 X) \quad (1)$$

where the  $W$  input layer maybe doing summation and summation just like matrix mutiplication does  $WX$ , and the  $\max$  function is a non-linear active function modifying the valuse like a neuron does before entering the next module:

$$W_2(\max(W_1 X)) \quad (2)$$

If now we don't have the active function then the formula will looks like:

$$W_2(W_1 X) \rightarrow \bar{W} X \quad (3)$$

which eventually all  $W_i$  can become a single module  $\bar{W}$

(b) For dictionary learning using sparse coding:

$$\min_{D, \alpha} \frac{1}{n} \sum_{i=1}^n \left( \frac{1}{2} \|x_i - D\alpha\|^2 + \lambda \|\alpha_i\|_1 \right) \quad (4)$$

Which is a joint minimization problem with respect to dictionary  $D$  and  $\alpha$

And for autoencoders:

$$\min_{W_{de}, W_{en}} \frac{1}{n} \sum_{i=1}^n \|W_{de} \sigma(W_{en} x_i) - x_i\|^2 \quad (5)$$

Where  $\sigma$  is some non-linear function (e.g. shrinkage). Similar to dictionary learning, autoencoders is also a joint optimization problem respect to encoder and decoder matrix  $W_{en}$ ,  $W_{de}$ . Indeed we can make the  $\alpha$  to become  $\sigma(W_{en} x)$  and this will make them very similar. However, there are two main factors making them different: One is autoencoder does not has the term for regulator, and therefore sparsity is not encouraged. Another one is autoencoder uses the model to find the code, while sparse coding approaching it by means of the optimizations.

## 2 Softmax regression gradient calculation

Given

$$\hat{y} = \sigma(Wx + b), \text{ where } x \in \mathbb{R}^d, W \in \mathbb{R}^{k \times d}, b \in \mathbb{R}^k \quad (6)$$

where  $d$  is the input dimension,  $k$  is the number of classes,  $\sigma$  is the softmax function:

$$\sigma(a)_i = \frac{\exp(a_i)}{\sum_j \exp(a_j)} \quad (7)$$

Which means a given input  $x$  will output  $y$  with probability of each class

21 **(a)** Derive  $\frac{\partial l}{\partial W_{ij}}$

22 If the given cross-entropy loss defined as followed:

$$l(y, \hat{y}) = - \sum_i y_i \log \hat{y}_i \quad (8)$$

23 As  $W_{ij}$  will affect the prediction of class  $i$  by multipling index  $j$  in  $x$ , therefore we can derive:

$$\frac{\partial l}{\partial W_{ij}} = \frac{\partial l}{\partial \hat{y}_i} \frac{\partial \hat{y}_i}{\partial W_{ij}} \quad (9)$$

24 where:

$$l(y, \hat{y}) = - \sum_i y_i \log \hat{y}_i = -(y_1 \log \hat{y}_1 + y_2 \log \hat{y}_2 + \dots + y_i \log \hat{y}_i + \dots) \quad (10)$$

25 and therefore

$$\frac{\partial l}{\partial \hat{y}_i} = \frac{-y_i}{\hat{y}_i} \quad (11)$$

26 And we can rewrite for only for  $\hat{y}_i$ :

$$\hat{y}_i = \frac{\exp(a_i)}{\sum_j \exp(a_j)} = \frac{\exp(a_i)}{C + \exp(a_i)}, \text{ where } C = \sum_{k \neq i} \exp(a_k) \quad (12)$$

27 Since

$$\frac{\partial \exp(a_i)}{\partial W_{ij}} = X_j \exp(a_i) \quad (13)$$

28 Therefore

$$\frac{\partial \hat{y}_i}{\partial W_{ij}} = X_j \hat{y}_i (1 - \hat{y}_i) \quad (14)$$

29 Finally, we will get the result of  $\frac{\partial l}{\partial W_{ij}}$ :

$$\frac{\partial l}{\partial W_{ij}} = \frac{\partial l}{\partial \hat{y}_i} \frac{\partial \hat{y}_i}{\partial W_{ij}} = -X_j y_i (1 - \hat{y}_i) \quad (15)$$

30 **(b)** What happen when  $y_{c_1} = 1, \hat{y}_{c_2} = 1, c_1 \neq c_2$

31 This means something like  $y = [1, 0, 0]^T$  and  $\hat{y} = [0, 0, 1]^T$ , and the predict is far different from true  
 32 label. This will cause the log part in loss (3) become negative infinity. We may not need to worry this  
 33 because before one of the class predicted close to 1 and everything else close to 0, it will generate a  
 34 great positive loss the the class that is miss-predicted trying to make the predict right to true label.

### 35 **3 Chain rule**

36 Without explicitly deriving the formula of  $f(x, y)$ , can we apply layers of functions to represent  
 37 function  $f$ , which is similar to build deep learning architecture.

$$\begin{aligned} f &= \frac{x^2 + \sigma(y)}{3x + y - \sigma(x)} = \frac{a}{b} \\ \implies \frac{\partial f}{\partial x} &= \frac{\partial a}{\partial x} \frac{1}{b} - \frac{a}{b^2} \frac{\partial b}{\partial x} \\ \implies \frac{\partial f}{\partial y} &= \frac{\partial a}{\partial y} \frac{1}{b} - \frac{a}{b^2} \frac{\partial b}{\partial y} \\ \implies \frac{\partial a}{\partial x} &= 2x \\ \implies \frac{\partial a}{\partial y} &= \sigma(y)(1 - \sigma(y)) \\ \implies \frac{\partial b}{\partial x} &= 3 - \sigma(x)(1 - \sigma(x)) \\ \implies \frac{\partial b}{\partial y} &= 1 \end{aligned} \quad (16)$$

38 **(b)** As  $x = 1$  and  $y = 0$ , then for each of value from the function listed above:

$$\begin{aligned}
a &= 1 + \sigma(0) = 1.5 \\
b &= 3 + 0 + \sigma(1) = 2.269 \\
\frac{\partial a}{\partial x} &= 2 \cdot 1 = 2 \\
\frac{\partial a}{\partial y} &= 0.5(1 - 0.5) = 0.25 \\
\frac{\partial b}{\partial x} &= 3 - 0.731(1 - 0.731) = 2.803 \\
\frac{\partial b}{\partial y} &= 1
\end{aligned} \tag{17}$$

39 Therefore, applying each of the gradient at  $(x, y) = (1, 0)$  to the chain rule, we will get:

$$\begin{aligned}
\frac{\partial f}{\partial x} &= \frac{\partial a}{\partial x} \frac{1}{b} - \frac{a}{b^2} \frac{\partial b}{\partial x} = 2 \cdot \frac{1}{2.269} - \frac{1.5}{(2.269)^2} \cdot 2.803 = 0.0647 \\
\frac{\partial f}{\partial y} &= \frac{\partial a}{\partial y} \frac{1}{b} - \frac{a}{b^2} \frac{\partial b}{\partial y} = 0.25 \cdot \frac{1}{2.269} - \frac{1.5}{(2.269)^2} \cdot 1 = -0.1811
\end{aligned} \tag{18}$$

#### 40 **4 Variants of pooling**

41 **(a)** The purpose of pooling is to progressively reducing the spatial size to reduce the amount  
42 of parameters and therefore also to control the issue of overfitting. There are many different  
43 variants of pooling for example max-pooling, average pooling, and fractional max-pooling, and  
44 they can be found in torch as function *SpatialMaxPooling*, *SpatialAveragePooling*, and  
45 *SpatialLPPooling*.

47 **(b)** For *SpatialMaxPooling* the definition is as followed:

$$x_{out} = \max(x_i^{(in)}) \quad \text{for signals in pool region} \tag{19}$$

48 For *SpatialAveragePooling* the definition is as followed:

$$x_{out} = \frac{1}{n} \sum_i^n x_i^{(in)} \quad \text{for signals in pool region} \tag{20}$$

49 For *SpatialLPPooling* the definition is as followed:

$$x_{out} = \frac{1}{n} \left( \sum_i^n (x_i^{(in)})^p \right)^{\frac{1}{p}} \quad \text{for signals in pool region} \tag{21}$$

50 **(c)** Max-pooling is very useful as it helps to eliminate non-maximal values and reduce the amount  
51 of parameter. However, if we just do max-pooling, the performance is limited due to its rapid  
52 reduction of spatial size, and the disjoint nature of the pooling region. Therefore, LP-pooling, which  
53 is an biologically inspired method, will be a moderate method that can reduce the spatial size as well  
54 as keeping the signal meaning in the pooling region.

#### 55 **5 Convolution**

56 **(a)** As it is using 3x3 kernel along x and y axis of input, which is 5 and 5 respectively. The output  
57 of this layer will be  $(5 - 3 + 1) \times (5 - 3 + 1)$  which is 3x3.

58

59 **(b)** Assuming the kernel operation is point-point multiplication and summation, then the output of  
60 this layer is:

$$61 \begin{pmatrix} 109 & 92 & 72 \\ 108 & 85 & 74 \\ 110 & 74 & 79 \end{pmatrix}$$

$$63 \text{ (c)} \begin{pmatrix} 4 & 7 & 10 & 6 & 3 \\ 9 & 17 & 25 & 16 & 8 \\ 11 & 23 & 34 & 23 & 11 \\ 7 & 16 & 24 & 17 & 8 \\ 2 & 6 & 9 & 7 & 3 \end{pmatrix}$$

## 65 **6 Optimization**

66 **(a)** say the encoder and decoder is defined as:

$$\begin{aligned} z &= W_1 x + b_1 \\ \tilde{x} &= W_2 z + b_2 \end{aligned} \quad (22)$$

67 And therefore the reconstruction loss  $J$  will be:

$$J(W_1, b_1, W_2, b_2) = (\tilde{x} - x)^2 = (W_2(W_1 x + b_1) + b_2 - x)^2 \quad (23)$$

68 **(b)** To have the gradient of reconstruction loss respective to the parameters, we take the derivative  
69 of each parameters:

$$\begin{aligned} \frac{\partial J}{\partial W_1} &= W_2 x \\ \frac{\partial J}{\partial W_2} &= W_1 x + b_1 \end{aligned} \quad (24)$$

70 **(c)** Say now we are at stage  $t$  and would like to compute  $W_1^{t+1}$  and  $W_2^{t+1}$ :

$$\begin{aligned} W_1^{t+1} &= W_1^t - \mu_1^t \frac{\partial J}{\partial W_1^t} = W_1^t - \mu_1^t (W_2 x) \\ W_2^{t+1} &= W_2^t - \mu_2^t \frac{\partial J}{\partial W_2^t} = W_2^t - \mu_2^t (W_1 x + b_1) \end{aligned} \quad (25)$$

71 where  $\mu_1^t$  and  $\mu_2^t$  are the step size at stage  $t$

72 **(d)** The updates during stochastic gradient descent usually involves Move-Forward and Correction  
73 stages and this oscillation may delay the efficiency of convergence, and therefore adding a momentum  
74 term may make the update toward the good direction as well as with the previous update history  
75 considered:

$$\begin{aligned} W_1^{t+1} &= W_1^t - \mu_1^t \frac{\partial J}{\partial W_1^t} + \Delta W_1^t \\ W_2^{t+1} &= W_2^t - \mu_2^t \frac{\partial J}{\partial W_2^t} + \Delta W_2^t \end{aligned} \quad (26)$$

## 76 **7 Top-k error**

77 For image classification, sometime the class is ambiguous, and the loss during is being modified to  
78 consider multiple label. The top-k error rate is the fraction of test images for which the correct label  
79 is not among the top-k labels considered most probable. The reason why ImageNet using both top-5  
80 and top-1 is due to sometimes only looking at top-1 error cannot be objective enough to evaluate the  
81 model because the image itself contains multi-label, and therefore evaluating top-5 error is important  
82 too.

## 83 8 t-SNE

84 (a) Suppose we have two-dimensional map that embedded with within a space with much higher  
 85 dimensionality, and if in this high dimension each data points are mutually equal distant, and the  
 86 mapping is not ablt toe faithfully performed if the convergence is simply just based on mutual distance.  
 87 Therefore, the crowding problem is defined as followed:

88 *The area of the two-dimensional map that is available to accommodate moderately distant datapoints*  
 89 *will not be nearly large enough compared with the area available to accommodate nearby datapoints.*

90 The approach of t-SNE to alleviate this problem is interesting. It converts distances into probability by  
 91 Gaussian distribution. Then , in low-dimension space, it uses a probability distribution that has much  
 92 heavier tails than Gaussian to conver distance to probability. After solving the joint optimization  
 93 problem, this model provides the conversion that data with small distances in high-dimension space  
 94 can get converted with much larger distances. Moreover, it eliminates the concerns of unwanted  
 95 attractive forces between dissimilar points.

96 There were attempts to solve the crowding problem, for example, by adding a small repulsion force  
 97 to all springs between data points. However, the optimization of this method is tedious, and also after  
 98 optimization it happend offently that two parts of cluster get seperated but then there is no force to  
 99 pull them back together.

100 (b) Derive  $\frac{\partial C}{\partial y_i}$

101 Let's define some variable for convenience:

$$\begin{aligned} d_{ij} &= ||y_i - y_j|| \\ Z &= \sum_{k \neq l} \exp(-d_{kl}^2) \end{aligned} \quad (27)$$

102 As we may notice that if we change  $y_i$  plus its symmetric property,  $d_{ij}$  and  $d_{ji}$  will be affected for  
 103  $\forall j$ . Therefore the gradient of C respect to  $y_i$  is given by:

$$\begin{aligned} \frac{\partial C}{\partial y_i} &= \sum_j \left( \frac{\partial C}{\partial d_{ij}} + \frac{\partial C}{\partial d_{ji}} \right) (y_i - y_j) \\ &= 2 \sum_j \frac{\partial C}{\partial d_{ij}} (y_i - y_j) \end{aligned} \quad (28)$$

104 Thus, the next question is to derive  $\frac{\partial C}{\partial d_{ij}}$ :

$$\begin{aligned} C &= \sum_i \sum_j (p_{ij} \log p_{ij} - p_{ij} \log q_{ij}) \\ \frac{\partial C}{\partial d_{ij}} &= - \sum_{k \neq l} p_{kl} \frac{\partial (\log q_{kl})}{\partial d_{ij}} = - \sum_{k \neq l} p_{kl} \frac{\partial (\log q_{kl} Z - \log Z)}{\partial d_{ij}} \\ &= - \sum_{k \neq l} p_{kl} \left( \frac{1}{q_{ij} Z} \frac{\partial (\exp(-d_{kl}^2))}{\partial d_{ij}} - \frac{1}{Z} \frac{\partial Z}{\partial d_{ij}} \right) \end{aligned} \quad (29)$$

105 Gradient  $\frac{\partial (\exp(-d_{kl}^2))}{\partial d_{ij}}$  is only nonzero when  $k = i$  and  $l = j$ . Therefore we can rewrite the formula  
 106 above to:

$$\begin{aligned} \frac{\partial C}{\partial d_{ij}} &= \frac{p_{ij}}{q_{ij} Z} (2 \exp(-d_{ij}^2)) - \sum_{k \neq l} \frac{2 \exp(-d_{ij}^2)}{Z} \\ &= 2p_{ij} - 2q_{ij} \end{aligned} \quad (30)$$

Therefore, the combine to the previous formula we will get:

$$\begin{aligned}\frac{\partial C}{\partial y_i} &= 2 \sum_j \frac{\partial C}{\partial d_{ij}} (y_i - y_j) \\ &= 4 \sum_j (p_{ij} - q_{ij}) (y_i - y_j)\end{aligned}\tag{31}$$

## 9 Proximal gradient descent

(a) Since Proximal operator is defined as:

$$\text{prox}_{h,t}(x) = \underset{z}{\operatorname{argmin}} \frac{1}{2} \|z - x\|_2^2 + th(z)\tag{32}$$

which the optimal condition is to have the gradient w.r.t  $z$  equal to 0:

$$0 \in z - x + t\partial h(z)\tag{33}$$

if function  $h(z) = \|z\|_1$  and  $z_i \neq 0$ , then:

$$\partial h(z) = \operatorname{sign}(z)\tag{34}$$

And therefore the optimal solution  $z^*$  will be:

$$z^* = x - t \cdot \operatorname{sign}(z^*)\tag{35}$$

Noted that if  $z_i^* < 0$ , then  $x_i < -t$ , and if  $z_i^* > 0$ , then  $x_i > t$ . This implies  $|x_i| > t$  and  $\operatorname{sign}(z_i^*) = \operatorname{sign}(x_i)$ , and we can rewrite formula to:

$$z_i^* = x_i - t \cdot \operatorname{sign}(x_i)\tag{36}$$

Then if the solution  $z_i^* = 0$ , the subgradient of l1-norm is in the interval of  $[-1, 1]$ , and we can write:

$$0 \in -x_i + t \cdot [-1, 1] \implies x_i \in [-t, t] \implies |x_i| \leq t\tag{37}$$

Therefore the solution of Proximal operator will be:

$$z_i^* = \begin{cases} 0 & \text{if } |x_i| \leq t \\ x_i - t \cdot \operatorname{sign}(x_i) & \text{if } |x_i| > t \end{cases}\tag{38}$$

which is

$$\text{prox}_{h,t}(x) = S_t(x) = (|x| - t)_+ \odot \operatorname{sign}(x) \quad (\text{element-wise})\tag{39}$$

which is a soft-threshold fuction with  $t$  as threshold value

(b) In the field of signal processing, the true signal usually will be blurred as followed:

$$Ax = b\tag{40}$$

where  $A$  is the blur operation,  $b$  is the known observed blurred-signal. The way to solve true signal  $x$  is called deblurring problem:

$$\min_x \{F(x) \equiv \frac{1}{2} \|b - Ax\|_2^2 + \lambda \|x\|_1\}\tag{41}$$

This is ISTA problem, and as we can see the first term is convex and differentiable, and the second term is convex and simple l1-norm function. Then the ISTA is become one example of proximal gradient descent

(c) From the definition of Proximal operator the optimal solution is where  $\frac{\partial \text{prox}_{h,t}}{\partial z} = 0$ , and therefore we will have:

$$0 \in z - x + t\partial h(z)\tag{42}$$

After we rewrite the function and replace  $z$  by  $u$  which is the optimal result from Proximal function:

$$\frac{x - u}{t} \in \partial h(u)\tag{43}$$

which means the calculated result from proximal function will be within the interval proportional to the subgradient of the simple-nonDerentiable function  $h(x)$

133 **(d)** From definition of Proximal operator, the optimal solution  $x_{k+1}$  will be:

$$x_{k+1} = \text{prox}_{h, \alpha_k}(x_k - \alpha_k \nabla g(x_k)) = x_k - \alpha_k \nabla g(x_k) - \alpha_k \partial h(x_{k+1}) \quad (44)$$

134 and from definition:

$$G_{\alpha_k}(x_k) = \frac{x_k - \text{prox}_{h, \alpha_k}(x_k - \alpha_k \nabla g(x_k))}{\alpha_k} \quad (45)$$

135 after rewrite:

$$x_k - \alpha_k \nabla g(x_k) - \alpha_k \partial h(x_{k+1}) = x_k - \alpha_k G_{\alpha_k}(x_k) \quad (46)$$

136 Therefore

$$G_{\alpha_k}(x_k) - \nabla g(x_k) \in \partial h(x_{k+1}) \quad (47)$$

137 which is because  $h$  is not differentiable and the result will within the range of subgradient of  $\partial h(x_{k+1})$