

Evaluating models fairly

MLP Tips

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Adapted from slides provided by Prof. Michael Mandel.

MLP design parameters

- Several parameters to choose when designing an MLP (best to evaluate empirically)
- Number of hidden layers
- Number of units in each hidden layer
- Activation function
- Error function

Optimization tricks

- For a given network, local minima of the cost function are possible
- Many tricks exist to try to find better local minima
 - Momentum: mix in gradient from step
 - Weight initialization: small random values
 - Stopping criterion: early stopping
 - Learning rate annealing: start with large, slowly shrink
 - Second order methods: use a separate for each parameter or pair of parameters based on local curvature
 - Randomization of training example order
 - Regularization, i.e., terms in $E(w)$ that only depend on w

Learning rate control: momentum

- To ease oscillating weights due to large η , some inertia (momentum) of weight update is added

$$\Delta w_{ji}(n) = \eta \delta_j y_i + \alpha \Delta w_{ji}(n-1), \quad 0 < \alpha < 1$$

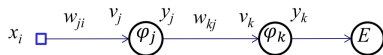
- In the downhill situation, $\Delta w_{ji}(n) \approx \frac{\eta}{1-\alpha} \delta_j y_i$
 - thus accelerating learning by a factor of $1/(1-\alpha)$
- In the oscillating situation, it smooths weight change, thus stabilizing oscillations

Input pre-processing

- Remove mean
 - Avoids extra update steps to learn it
- Divide by standard deviation
 - Or whiten by multiplying by the inverse of square root of the covariance matrix
 - Make dimensions commensurate
 - Scales curvature of error surface to be less canyon-like

Weight initialization

- Consider a network with one hidden layer and a single output neuron
- What happens if we initialize all weights to 0?

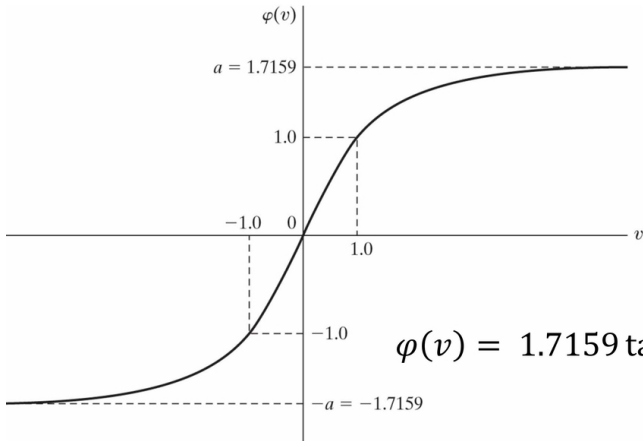


$$y_k = \varphi_k \left(\sum_j w_{kj} \varphi_j \left(\sum_i w_{ji} x_i \right)_j \right)$$
$$\frac{\partial}{\partial w_{ki}} E(\mathbf{w}) = -e_k \varphi'(v_k) y_j$$
$$\frac{\partial}{\partial w_{ji}} E(\mathbf{w}) = -e_j \varphi'(v_j) x_i$$

Weight initialization

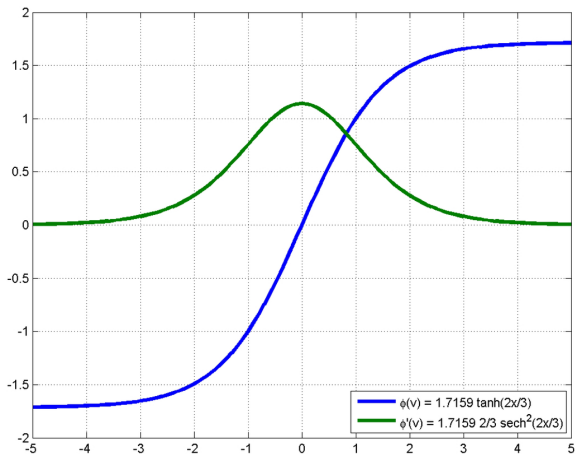
- Break symmetry by initializing with random values
- If inputs are normalized, they are uncorrelated, with zero-mean, and unit-variance, and all weights are initialized to 0,
- Then, every neuron in the network computes the same output, and also computes the same gradients during backpropagation and undergoes the exact same parameter updates. The network never converges.
- Break symmetry accelerate convergence.

Hyperbolic tangent function



$$\varphi(v) = 1.7159 \tanh\left(\frac{2}{3}v\right)$$

Hyperbolic tangent function



Weight initialization

$$\begin{aligned}\sigma_{y_i}^2 &= E_x\{y_i^2\} = E_x\left\{\varphi^2\left(\sum_j w_{ij}x_j\right)\right\} \\ &\approx E_x\left\{\left(\sum_j w_{ij}x_j\right)^2\right\} \approx \sum_j w_{ij}^2 E_x\{x_j^2\} \\ &= \sum_{j=1}^m w_{ij}^2\end{aligned}$$

- So in order to make $\sigma_{y_i}^2 = 1$
 - Initialize w_{ij} randomly with $\sigma_w^2 = \frac{1}{m}$

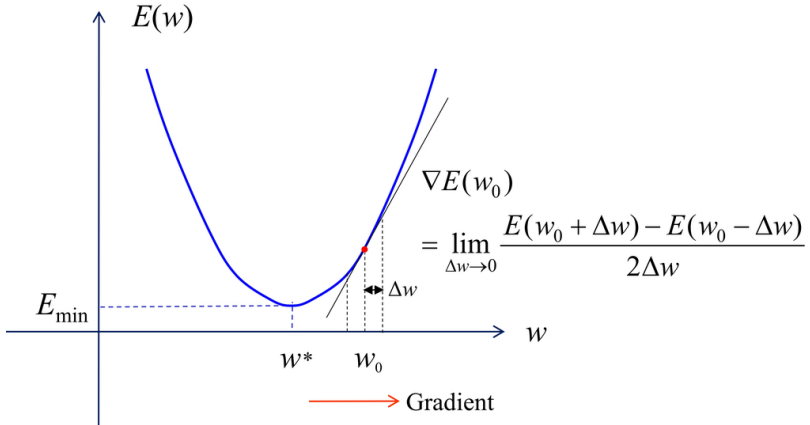
Debugging: Gradient checking

- Is your backpropagation code working properly?
 - *i.e.*, is it computing the right gradient?
- Backpropagation computes

$$\nabla_{\mathbf{w}} E(\mathbf{x}_p; \mathbf{w}) = \left[\frac{\partial E}{\partial \mathbf{w}_{111}}, \frac{\partial E}{\partial \mathbf{w}_{121}}, \dots, \frac{\partial E}{\partial \mathbf{w}_{NML}} \right]$$

- where $w_{i_1 i_2 l}$ is the weight in layer l connecting neurons i_1 and i_2
- Compute the gradient **numerically** and compare

Recall: Gradient illustration



Debugging: Gradient checking

- One-sided numerical gradient:

$$\frac{\partial E}{\partial w_{i_1 i_2 l}} \approx \frac{1}{\delta} \left(E(\mathbf{x}_p; \mathbf{w} + \delta \mathbf{1}_{i_1 i_2 l}) - E(\mathbf{x}_p; \mathbf{w}) \right)$$

- where $\mathbf{1}_{i_1 i_2 l}$ is a vector that is 1 at entry $i_1 i_2 l$ and 0 everywhere else and δ is a “small” constant
- Two-sided numerical gradient:

$$\frac{1}{2\delta} \left(E(\mathbf{x}_p; \mathbf{w} + \delta \mathbf{1}_{i_1 i_2 l}) - E(\mathbf{x}_p; \mathbf{w} - \delta \mathbf{1}_{i_1 i_2 l}) \right)$$

- More accurate approximation
- But requires twice as many evaluations of $E(\mathbf{x}_p; \mathbf{w})$

Debugging: Gradient checking

- Complexity of backpropagation
 - 1 forward pass ($O(1)$ multiply and add per weight)
 - 1 backward pass ($O(1)$ multiply and add per weight)
- Complexity of numerical gradient
 - One-sided: 1 forward pass *per network weight*
 - So $W + 1$ forward passes total
 - Two-sided: 2 forward passes *per network weight*
- So numerical gradient is good for checking correctness of backpropagation
 - But very slow to use in training, especially for large W

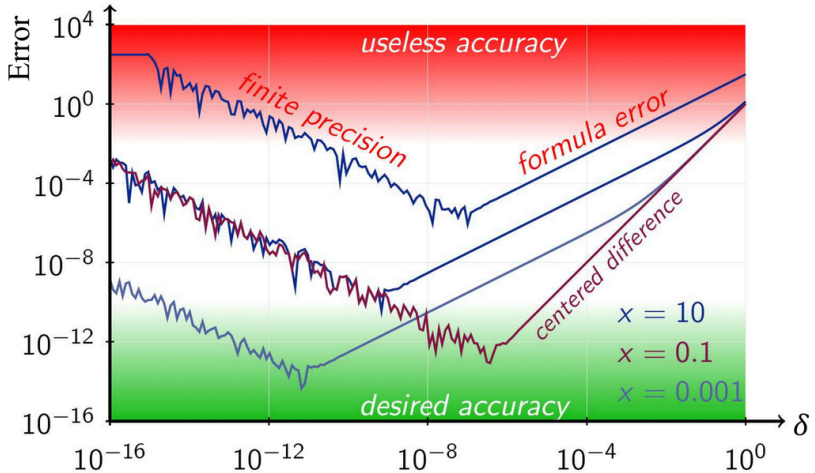
Gradient checking procedure

- Select an example data point, \mathbf{x}_p , initialize \mathbf{w}
- Compute the gradient of $E(\mathbf{x}_p; \mathbf{w})$ using backprop
 - Gives a vector of derivatives, one for each weight in the network
- Compute the gradient numerically
 - Evaluate $E(\mathbf{x}_p; \mathbf{w})$
 - Loop over each weight in the network
 - Evaluate $E(\mathbf{x}_p; \mathbf{w} + \delta \mathbf{1}_{i_1 i_2 l})$, compute partial derivative
- If they are not the same, look for patterns as a function of $i_1 i_2 l$, etc

How to select δ ?

- δ too big means derivative might be different at $E(\mathbf{x}_p; \mathbf{w} + \delta \mathbf{1}_{i_1 i_2 l})$ and $E(\mathbf{x}_p; \mathbf{w})$
 - Leading to a bad estimate using the above formulas
- δ too small runs into numerical issues
 - Need to be aware of limitations of floating point math
 - For δ too small, $1 + \delta = 1$
 - This might be around $1\text{e-}16$, depending on the data type (e.g., float, double)
 - So $\delta = 1\text{e-}8$ might be reasonable

How to select δ ?



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Thank you!