Applied Machine Learning

Gradient Computation & Automatic Differentiation

Oumar Kaba



Learning objectives

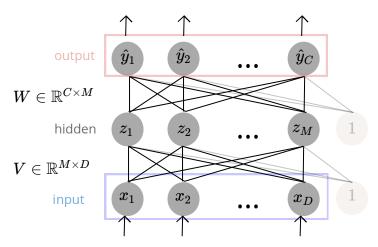
using the chain rule to calculate the gradients automatic differentiation

- forward mode
- reverse mode (backpropagation)

Landscape of the cost function

model two layer MLP

$$f(x;W,V) = g(Wh(Vx))$$



for simplicity we drop the bias terms

objective
$$\min_{W,V} \sum_n L(y^{(n)}, f(x^{(n)}; W, V))$$
 loss function depends on the task

this is a non-convex optimization problem



https://losslandscape.com/gallery/

Landscape of the cost function

model two layer MLP

$$f(x;W,V) = gig(Wh(Vx)ig)$$

general beliefs

supported by empirical and theoretical results in a special settings

many more saddle points than local minima number of local minima increases for lower costs therefore most local optima are close to global optima

strategy

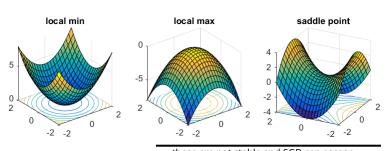
use gradient descent methods (covered earlier in the course)

objective $\min_{W,V} \sum_n L(y^{(n)}, f(x^{(n)}; W, V))$

this is a non-convex optimization problem



many critical points (points where gradient is zero)



these are not stable and SGD can escape

Jacobian matrix

Examples

 $f: \mathbb{R} o \mathbb{R}$ - we have the derivative $rac{d}{dw} f(w) \in \mathbb{R}$

 $f:\mathbb{R}^D o\mathbb{R}$ gradient is the vector of all partial derivatives

$$abla_w f(w) = [rac{\partial}{\partial w_1} f(w), \ldots, rac{\partial}{\partial w_D} f(w)]^ op \in \mathbb{R}^D$$

 $f:\mathbb{R}^D o\mathbb{R}^M$ the **Jacobian matrix** of all partial derivatives

$$rac{\partial}{\partial w_1}f(w)$$
 $\nabla_w f_1(w)$ $\left[egin{array}{c} rac{\partial f_1(w)}{\partial w_1}, & \dots, & rac{\partial f_1(w)}{\partial w_D} \end{array}
ight]$ $J=\left[egin{array}{c} dots & \ddots & dots & d$

$$f(x) = x^2 \Rightarrow \frac{d}{dw}f(w) = 2x$$

$$f(x,y) = x^2 + y^2 \Rightarrow
onumber
abla_{w=[x,y]} f(w) = [2x,2y]$$

$$f(x,y) = \begin{bmatrix} x^2, y^2, 2xy \end{bmatrix} \Rightarrow$$

$$J = \begin{bmatrix} 2x, & 0 \\ 0, & 2y \\ 2y, & 2x \end{bmatrix}$$

for all three case we may simply write $\frac{\partial}{\partial w}f(w)$, where M,D will be clear from the context what if W is a matrix? we assume it is reshaped into a vector for these calculations

Chain rule

for
$$f: x \mapsto z$$
 and $h: z \mapsto y$ where $x,y,z \in \mathbb{R}$

$$\frac{dy}{dx} = \frac{dy}{dz} \frac{dz}{dx}$$

$$| speed of change in z as we change x$$

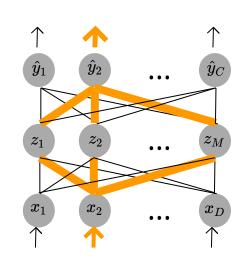
$$| speed of change in y as we change z$$

$$| speed of change in y as we change x$$

more generally $x \in \mathbb{R}^D, z \in \mathbb{R}^M, y \in \mathbb{R}^C$

$$rac{\partial y}{\partial x} = rac{\partial y}{\partial z} rac{\partial z}{\partial x}$$
 in matrix form
 $C \, x \, D \, Jacobian$ $M \, x \, D \, Jacobian$ $C \, x \, M \, Jacobian$

$$rac{\partial y_c}{\partial x_d} = \sum_{m=1}^{M} rac{\partial y_c}{\partial z_m} rac{\partial z_m}{\partial x_d}$$



Training a two layer network

model
$$\hat{y} = gig(W\,h(V\,x)ig)$$

Cost function we want to minimize

$$J(W,V) = \sum_n L(y^{(n)}, g(W h(V x^{(n)}))$$

need gradient wrt W and V: $rac{\partial}{\partial W}J, \; rac{\partial}{\partial V}J$

output \hat{y}_1 \hat{y}_2 ... \hat{y}_C W hidden units z_1 z_2 ... z_M 1 v input v ... v 1

for simplicity we drop the bias terms

simpler to write this for one instance (n)

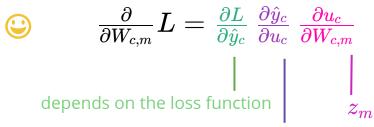
so we will calculate $\ \frac{\partial}{\partial W}L, \ \frac{\partial}{\partial V}L$ and recover

$$rac{\partial}{\partial W} J = \sum_{n=1}^N rac{\partial}{\partial W} L(y^{(n)}, \hat{y}^{(n)})$$
 and $rac{\partial}{\partial V} J = \sum_{n=1}^N rac{\partial}{\partial V} L(y^{(n)}, \hat{y}^{(n)})$



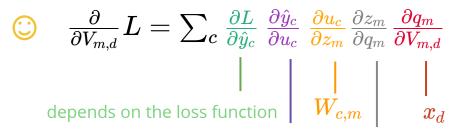


using the chain rule



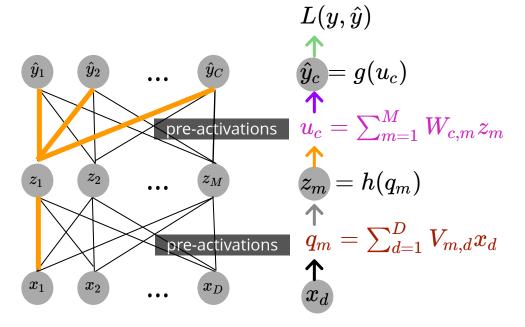
depends on the activation function

similarly for V



depends on the activation function

depends on the middle layer activation



using the chain rule

$$rac{\partial}{\partial W_{c,m}} L = rac{\partial L}{\partial \hat{y}_c} rac{\partial \hat{y}_c}{\partial u_c} rac{\partial u_c}{\partial W_{c,m}}$$
 depends on the loss function depends on the activation function z_m $C=1$ $L(y,\hat{y}) = rac{1}{2}||y-\hat{y}||_2^2$ $\hat{y}=g(u)=u$

combining the three terms above

$$rac{\partial}{\partial W_m}L=(\hat{y}-y)z_m$$
 we have seen this in linear regression lecture!

$$egin{aligned} L(y,\hat{y}) \ \hat{y}_c &= g(u_c) \ v_c &= \sum_{m=1}^M W_{c,m} z_m \ z_m &= h(q_m) \ q_m &= \sum_{d=1}^D V_{m,d} x_d \ x_d \end{aligned}$$

more generally:

$$rac{\partial}{\partial W_{c,m}}L=(\hat{y}_c-y_c)z_m$$
 ,

using the chain rule



$$rac{\partial}{\partial W_{c,m}}L=rac{\partial L}{\partial \hat{y}_c}rac{\partial \hat{y}_c}{\partial u_c}rac{\partial u_c}{\partial W_{c,m}}$$
 depends on the loss function depends on the activation function

binary classification

$$egin{align} L(y,\hat{y}) &= -y\log\hat{y} - (1-y)\log(1-\hat{y}) & rac{\partial\hat{y}}{\partial u} &= \hat{y}(1-\hat{y}) \ rac{\partial}{\partial\hat{y}}L(y,\hat{y}) &= -rac{y}{\hat{y}} + rac{(1-y)}{(1-\hat{y})} \ \end{pmatrix} \end{array}$$

combining the three terms above

$$rac{\partial}{\partial W_m} L = (\hat{y} - y) z_m$$

 $L(y,\hat{y})$ $egin{aligned} \hat{y}_c &= g(u_c) \ v_c &= \sum_{m=1}^M W_{c,m} z_m \end{aligned}$ $egin{aligned} oldsymbol{\uparrow} \ oldsymbol{z}_m &= h(q_m) \ oldsymbol{\uparrow} \ oldsymbol{q}_m &= \sum_{d=1}^D V_{m,d} x_d \end{aligned}$

using the chain rule



$$rac{\partial}{\partial W_{c,m}}L=\sum_{k=1}^Crac{\partial L}{\partial \hat{y}_k}rac{\partial \hat{y}_k}{\partial u_c}rac{\partial u_c}{\partial W_{c,m}}$$
 depends on the loss function depends on the activation function z_m

multiclass classification

C is the number of classes

$$L(y,\hat{y}) = -\sum_c y_c \log \hat{y}_c \;\; \hat{y} = g(u) = ext{softmax}(u) \;\; ext{softmax takes a vector and produces a vector} \ rac{\partial}{\partial \hat{y}_k} L = -rac{y_k}{\hat{y}_k} \;\;\; \hat{y}_k = rac{e^{u_k}}{\sum_i e^{u_i}} \;\; ext{need to calculate the Jacobian} \;\; rac{\partial}{\partial u_c} \hat{y}_k = egin{cases} \hat{y}_k (1 - \hat{y}_k) & k = c \\ -\hat{y}_c \hat{y}_k & k \neq c \end{cases}$$

combining the three terms above

$$rac{\partial}{\partial W_{c,m}}L=(\hat{y}_c-y_c)z_m$$

 $L(y,\hat{y})$

 $egin{aligned} oldsymbol{\hat{y}_c} &= g(u_c) \ oldsymbol{\uparrow} \ u_c &= \sum_{m=1}^M W_{c,m} z_m \end{aligned}$

 $egin{aligned} oldsymbol{\uparrow} & oldsymbol{\uparrow} \ oldsymbol{z}_m = h(q_m) \ oldsymbol{\uparrow} \ q_m = \sum_{d=1}^D V_{m,d} x_d \end{aligned}$

gradient wrt V:

we already did this part



$$rac{\partial}{\partial V_{m,d}} L = \sum_{c} \left[egin{array}{ccccc} rac{\partial L}{\partial \hat{y}_{c}} & rac{\partial \hat{y}_{c}}{\partial u_{c}} & rac{\partial u_{c}}{\partial z_{m}} & rac{\partial z_{m}}{\partial q_{m}} & rac{\partial q_{m}}{\partial V_{m,d}}
ight] \ W_{c,m} & \left| egin{array}{c} I & x_{d} \end{array}
ight|$$

depends on the middle layer activation

logistic function
$$\sigma(q_m)(1-\sigma(q_m))$$
 hyperbolic tan. $1- anh(q_m)^2$ ReLU $egin{cases} 0 & q_m \leq 0 \ 1 & q_m > 0 \end{cases}$

example

logistic sigmoid

$$egin{aligned} rac{\partial}{\partial V_{m,d}} L &= \sum_c (\hat{y}_c - y_c) W_{c,m} \sigma(q_m) (1 - \sigma(q_m)) oldsymbol{x_d} \ &= \sum_c (\hat{y}_c - y_c) W_{c,m} z_m (1 - z_m) oldsymbol{x_d} \end{aligned} \qquad \Rightarrow rac{\partial}{\partial V_{m,d}} J = \sum_n \sum_c (\hat{y}_c^{(n)} - y_c^{(n)}) W_{c,m} z_m^{(n)} (1 - z_m^{(n)}) oldsymbol{x_d^{(n)}} \ \end{aligned}$$

 $L(y,\hat{y})$

 $egin{aligned} \hat{y_c} &= g(u_c) \ v_c &= \sum_{m=1}^M W_{c,m} z_m \end{aligned}$

 $egin{aligned} oldsymbol{\uparrow} \ oldsymbol{z}_m &= h(q_m) \ oldsymbol{\uparrow} \ q_m &= \sum_{d=1}^D V_{m,d} x_d \end{aligned}$

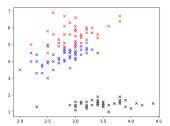
a common pattern

$$\stackrel{\partial}{\partial W_{c,m}} L = \left[egin{array}{ccc} rac{\partial L}{\partial \hat{y}_c} & rac{\partial y_c}{\partial u_c} & rac{\partial u_c}{\partial W_{c,m}}
ight] \ & ext{error from above} & rac{\partial L}{\partial u_c} & ext{input from below} & z_m \end{array}$$

$$\stackrel{\partial}{\partial V_{m,d}} L = \sum_{c} rac{\partial L}{\partial \hat{y}_{c}} rac{\partial \hat{y}_{c}}{\partial u_{c}} rac{\partial u_{c}}{\partial z_{m}} rac{\partial z_{m}}{\partial q_{m}} rac{\partial q_{m}}{\partial V_{m,d}}$$
 error from above $rac{\partial L}{\partial q_{m}}$ input from below x_{d} .

$$egin{aligned} L(y,\hat{y}) \ \hat{m{y}_c} &= g(u_c) \ m{u}_c &= \sum_{m=1}^M W_{c,m} z_m \ m{z}_m &= h(q_m) \ m{q}_m &= \sum_{d=1}^D V_{m,d} x_d \ m{\uparrow} \ m{x}_d \end{aligned}$$

Example: classification



Iris dataset (D=2 features + 1 bias)

M = 16 hidden units

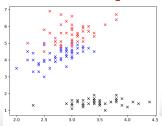
C=3 classes

cost is softmax-cross-entropy

$$L(y,\hat{y})$$
 $\hat{y} = \operatorname{softmax}(u)$
 $u_c = \sum_{m=1}^M W_{c,m} z_m$
 $z_m = \sigma(q_m)$
 $q_m = \sum_{d=1}^D V_{m,d} x_d$
 x_d

$$J = -\sum_{n=1}^{N} y^{(n)}^{ op} u^{(n)} + \log \sum_{c} e^{u_{c}^{(n)}}$$

Example: classification



10

11

12 13

```
Iris dataset (D=2 features + 1 bias)
```

M = 16 hidden units

C=3 classes

 $dy = yh - y \#N \times C$

return dw, dv

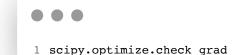
dw= np.dot(z.T, dy)/N #M x C

dz = np.dot(dy, w.T) #N x M

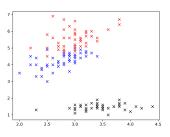
dv = np.dot(x.T, dz * z * (1 - z))/N #D x M

$$L(y, \hat{y})$$
 $\hat{y} = \operatorname{softmax}(u)$
 $u_c = \sum_{m=1}^{M} W_{c,m} z_m$
 $z_m = \sigma(q_m)$
 $q_m = \sum_{d=1}^{D} V_{m,d} x_d$
 x_d

check your gradient function using **finite difference** approximation that uses the *cost function*



Example: classification

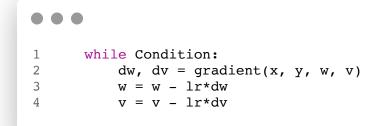


Iris dataset (D=2 features + 1 bias)

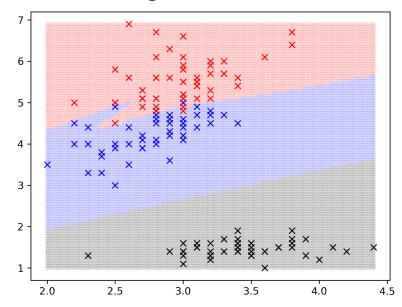
M = 16 hidden units

C=3 classes

using GD for optimization



the resulting decision boundaries



Automating gradient computation

gradient computation is tedious and mechanical. Can we automate it?

using numerical differentiation?

approximates partial derivatives using finite difference $\frac{\partial f}{\partial w} \approx \frac{f(w+\epsilon)-f(w)}{\epsilon}$ needs multiple forward passes (for each input output pair) can be slow and inaccurate useful for black-box cost functions or checking the correctness of gradient functions

symbolic differentiation: symbolic calculation of derivatives

does not identify the computational procedure and reuse of values

automatic / algorithmic differentiation is what we want

write code that calculates various functions, *e.g., the cost function* automatically produce (partial) derivatives *e.g., gradients used in learning*

Automatic differentiation

idea use the chain rule + derivative of simple operations $*, \sin, \frac{1}{x}$...

use a computational graph as a data structure (for storing the result of computation)

step 1 break down to atomic operations

$$L=rac{1}{2}(wx-y)^2$$
 $a_1=w$ $a_2=x$

- step 2 build a graph with operations as internal nodes and input variables as leaf nodes
- step 3 there are two ways to use the computational graph to calculate derivatives

forward mode: start from the leafs and propagate derivatives upward

reverse mode:

- 1. first in a bottom-up (forward) pass calculate the values $\,a_1,\ldots,a_4\,$
- 2. in a top-down (backward) pass calculate the derivatives

this second procedure is called **backpropagation** when applied to neural networks

Forward mode

suppose we want the derivative
$$\ rac{\partial y_1}{\partial w_1} \$$
 where $\ egin{cases} y_1 = \sin(w_1x + w_0) \ y_2 = \cos(w_1x + w_0) \end{cases}$

we can calculate both y_1, y_2 and derivatives $\frac{\partial y_1}{\partial w_1}$ $\frac{\partial y_2}{\partial w_1}$ in a single forward pass

evaluation

$$a_2 = w_1$$

$$a_3 = x$$

$$w_1x \hspace{1cm} a_4=a_2 imes a_3 \hspace{1cm} \dot{a_4}=a_2 imes \dot{a_3}+\dot{a_2} imes a_3$$

$$w_1 x + w_0$$
 $a_5 = a_4 + a_1$ $a_5 = a_4 + a_1$

$$a_0=\sin(w_1x+w_0)$$
 $a_0=\sin(a_0)$

$$a_2=\cos(w_1x+w_0)$$
 $a_7=\cos(a_5)$

partial derivatives

$$egin{array}{l} \dot{a_1}=0 \ \dot{a_2}=1 \ \dot{a_3}=0 \end{array}
ight\} egin{array}{l} \mathsf{we} \ \mathsf{thi} \end{array}$$

$$\dot{a}_4 = a_2 \times \dot{a}_3 + \dot{a}_2 \times a_3 \qquad x$$

$$i_5 = \dot{a_4} + \dot{a_1} \qquad \qquad x$$

$$-x\sin(w_1x+w_0)=rac{\partial y_2}{\partial w_1}$$

Forward mode: computational graph

suppose we want the derivative $\frac{\partial y_1}{\partial w_1}$ where $\begin{cases} y_1 = \sin(w_1 x + w_0) \\ y_2 = \cos(w_1 x + w_0) \end{cases}$

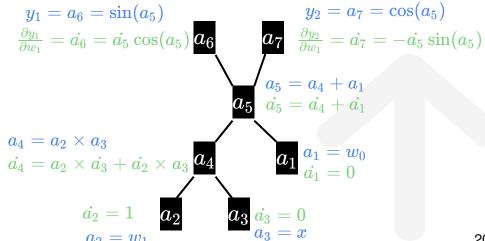
we can represent this computation using a graph once the nodes up stream calculate their values and derivatives we may discard a node

• e.g., once $a_5, \dot{a_5}$ are obtained we can discard the values and partial derivatives for $a_4, \dot{a_4}, a_1, \dot{a_1}$

evaluation $a_1 = w_0$ $a_2 = w_1$ $a_3 = x$ $a_4 = a_2 \times a_3$ $y_1 = a_6 = \sin(a_5)$ $a_6 = a_5 \cos(a_5)$

partial derivatives

$$egin{array}{lll} a_1 = w_0 & \dot{a_1} = 0 \ a_2 = w_1 & \dot{a_2} = 1 \ a_3 = x & \dot{a_3} = 0 \ a_4 = a_2 imes a_3 & \dot{a_4} = a_2 imes \dot{a_3} + \dot{a_2} imes a_3 \ a_5 = a_4 + a_1 & \dot{a_5} = \dot{a_4} + \dot{a_1} \ y_1 = a_6 = \sin(a_5) & \dot{a_6} = \dot{a_5}\cos(a_5) \ y_2 = a_7 = \cos(a_5) & \dot{a_7} = -\dot{a_5}\sin(a_5) \ \end{array}$$



Reverse mode

suppose we want the derivative $\ \, rac{\partial y_2}{\partial w_1} \,$ where $\ \, y_2 = \cos(w_1 x + w_0) \,$

 w_1x

2) partial derivatives

first do a forward pass for evaluation

1) evaluation

 $a_1 = w_0$

then use these values to calculate partial derivatives in a backward pass

 $a_2 = w_1$

 $a_3 = x$

 $a_4 = a_2 \times a_3$

 $rac{\partial y_2}{\partial a_7}=1$ $ar{a_7}=1$ this means $ar{\Box}=rac{\partial y_2}{\partial \Box}$ $rac{\partial y_2}{\partial a_8}=0$

 $\frac{\partial y_2}{\partial w_1} = \frac{\partial y_2}{\partial a_2} = \frac{\partial y_2}{\partial a_4} = \frac{\partial a_4}{\partial a_2} = -x \sin(w_1 x + w_0)$ $\bar{a_2} = \bar{a_4} a_3$ $\frac{\partial y_2}{\partial w_0} = \frac{\partial y_2}{\partial a_1} = \frac{\partial y_2}{\partial a_2} \frac{\partial a_5}{\partial a_1} = -\sin(w_1 x + w_0)$ $\bar{a_1} = \bar{a_5}$

we get all partial derivatives $\frac{\partial y_2}{\partial \Box}$ in one backward pass

21

Reverse mode: computational graph

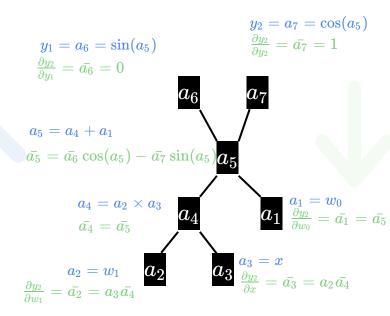
suppose we want the derivative $\ \, rac{\partial y_2}{\partial w_1} \,$ where $\ \, y_2 = \cos(w_1 x + w_0) \,$

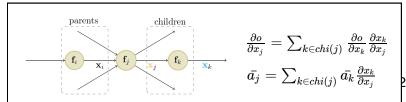
we can represent this computation using a graph

- 1. in a forward pass we do evaluation and **keep the values**
- 2. use these values in the backward pass to get partial derivatives

1) evaluation

$$a_1 = w_0$$
 $a_2 = w_1$ $a_{\bar{7}} = 1$ $a_{\bar{6}} = 0$ $a_{\bar{3}} = x$ $a_{\bar{6}} = 0$ $a_{\bar{5}} = \bar{a}_{\bar{6}} \cos(a_{\bar{5}}) - \bar{a}_{\bar{7}} \sin(a_{\bar{5}})$ $a_{\bar{5}} = a_{\bar{4}} + a_{\bar{1}}$ $a_{\bar{5}} = a_{\bar{6}} \cos(a_{\bar{5}}) - \bar{a}_{\bar{7}} \sin(a_{\bar{5}})$ $a_{\bar{5}} = a_{\bar{4}} + a_{\bar{1}}$ $a_{\bar{4}} = \bar{a}_{\bar{5}}$ $a_{\bar{3}} = a_{\bar{2}} \bar{a}_{\bar{4}}$ $a_{\bar{2}} = a_{\bar{3}} \bar{a}_{\bar{4}}$ $a_{\bar{2}} = a_{\bar{3}} \bar{a}_{\bar{4}}$ $a_{\bar{1}} = \bar{a}_{\bar{5}}$

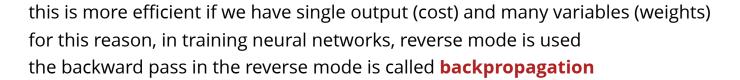




Forward vs Reverse mode

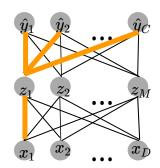
forward mode is more natural, easier to implement and requires less memory a single **forward** pass calculates $\frac{\partial y_1}{\partial w}, \dots, \frac{\partial y_c}{\partial w}$

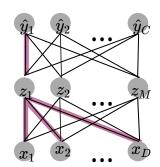
however, **reverse** mode is more efficient in calculating gradient $\nabla_w y = [rac{\partial y}{\partial w_1}, \dots, rac{\partial y}{\partial w_D}]^ op$



many machine learning software implement autodiff:

- autograd (extends numpy)
- pytorch
- tensorflow





Improving optimization in deep learning

Initialization of parameters:

- random initialization (uniform or Gaussian) with small variance
 - break the symmetry of hidden units
- small positive values for bias (so that input to ReLU is >0)

models that are simpler to optimize:

this block is correcting for the residual error in the predictions of the previous layers

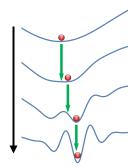


• using **skip-connection**
$$x^{\{\ell+l\}} = \mathrm{ReLU}(W^{\{\ell+l\}}\mathrm{ReLU}(\dots\mathrm{ReLU}(W^{\{\ell\}}x^{\{\ell\}})\dots) + x^{\{\ell\}})$$

using batch-normalization (next)

Pretrain a (simpler) model on a (simpler) task and

fine-tune on a more difficult target setting (has many forms)



continuation methods in optimization

- gradually increase the difficulty of the optimization problem
- good initialization for the next iteration

curriculum learning (similar idea)

- increase the number of "difficult" examples over time
- similar to the way humans learn

weight layer

weight layer

∓relu

 \mathbf{x}

identity

Batch Normalization

original motivation

- gradient descent: parameters in all layers are updated
- distribution of inputs to layer ℓ changes
- each layer has to re-adjust
- inefficient for very deep networks

activation for the instance (n) at layer ℓ

normalize the input to each unit (m) of a layer ℓ idea

alternatively: apply the batch-norm to $W^{\{\ell\}}x^{\{\ell\}}$

each unit is unnecessarily constrained to have zero-mean and std=1 (we only need to fix the distribution)

introduce learnable parameters
$$\operatorname{ReLU}(\gamma^{\{\ell\}} \operatorname{BN}(W^{\{\ell\}} x^{\{\ell\}}) + eta^{\{\ell\}})$$

- mean and std per unit is calculated for the minibatch during the forward pass
- we backpropagate through this normalization
- at test time use the mean and std. from the whole training set
- BN regularizes the model

recent observations the change in distribution of activations is not a big issue empirically BN works so well because it makes the loss function smooth

Summary

optimization landscape in neural networks is special and not yet fully understood

- exponentially many local optima and saddle points
- most local minima are good
- calculate the gradients using backpropagation

automatic differentiation

- simplifies gradient calculation for complex models
- gradient descent becomes simpler to use
- ullet forward mode is useful for calculating the jacobian of $\,f:\mathbb{R}^Q o\mathbb{R}^P$ when $\,P\geq Q\,$
- reverse mode can be more efficient when Q>P
 - backpropagation is reverse mode autodiff.

Better optimization in deep learning:

- better initialization
- models that are easier to optimize (using skip-connection, batch-norm, ReLU)
- pre-training and curriculum learning