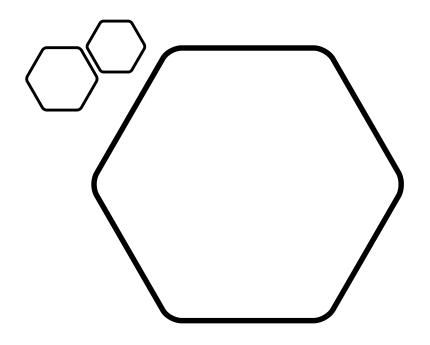
Lecture 7

Cs231n



SGD

Optimization

```
# Vanilla Gradient Descent

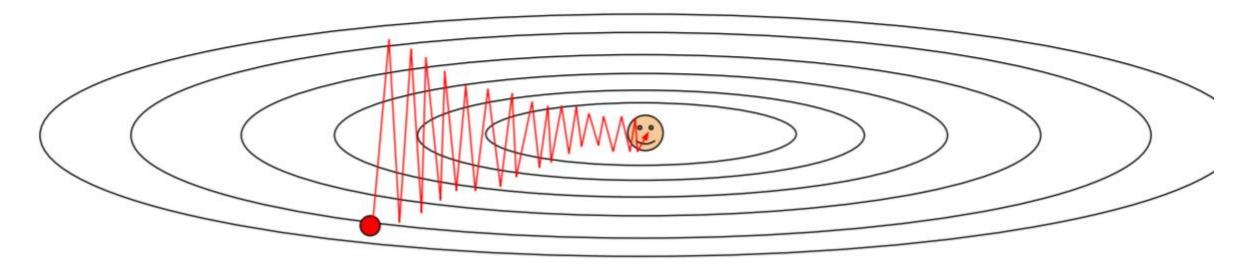
while True:
    weights_grad = evaluate_gradient(loss_fun, data, weights)
    weights += - step_size * weights_grad # perform parameter update
```

W_2

W_1

Problem with SGD – loss function affects a lot

• IF) 한 방향의 가중치에 대해서만 잘 영향받는 descent?

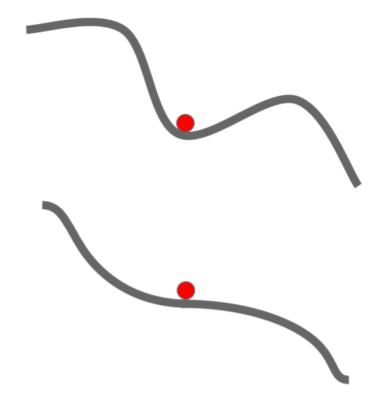


• 지그재그로 움직이는 문제 발생

Problem with SGD

• IF) local minima 나 saddle point(안장점) 에 안착할 경우 더 좋은 값을 안찾으러감.

 더 큰 문제는 high dimension 일수록 더 많은 local minima, saddle point 가 있음.

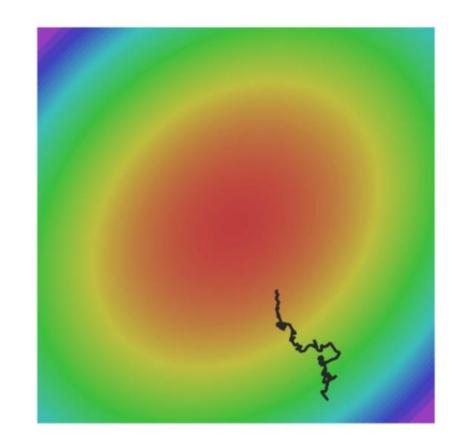


Problem with SGD

• SGD 는 배치단위 → 노이즈가 발생할 확률이 올라감. → 시간이 많이걸리게됨.

$$L(W) = \frac{1}{N} \sum_{i=1}^{N} L_i(x_i, y_i, W)$$

$$\nabla_W L(W) = \frac{1}{N} \sum_{i=1}^{N} \nabla_W L_i(x_i, y_i, W)$$



이 모든 문제는 full batch에서도 발생함.

몇가지 해결방법

- Momentum
 - -Nesterov Momentum
- AdaGrad
- RMSProp
- Adam

Momentum

- 운동량의 개념을 도입. "속도", "마찰력"의 개념을 생각
- 모든 문제가 조금은 완화됨.
- local minima, saddle point에서도 속도를 가지고있다고 생각하면 멈추지 않고 계속감
- 가속도의 개념으로 poor conditioning 을 해결가능(지그재그로 가지않는 지점은 가속) sgd sgD+Momentum
- 모멘텀이 노이즈를 평균화하여 완화

 $x_{t+1} = x_t - \alpha \nabla f(x_t)$

```
x_{t+1} = x_t - \alpha v_{t+1}  \text{vx = 0}  while True:
```

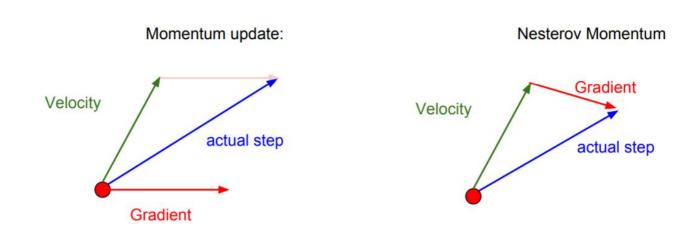
 $v_{t+1} = \rho v_t + \nabla f(x_t)$

```
dx = compute_gradient(x)
vx = rho * vx + dx
x += learning_rate * vx
```

- Build up "velocity" as a running mean of gradients
- Rho gives "friction"; typically rho=0.9 or 0.99

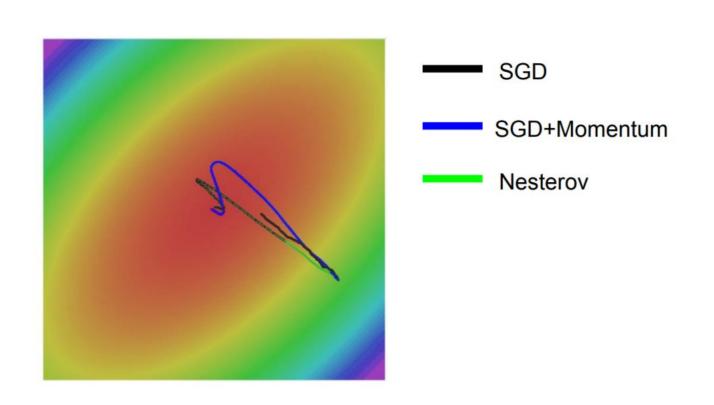
Nesterov Momentum

- Actural update는 gradient와 속 도의 가중평균으로 구해짐
- 기본 아이디어는 같으나 순서를 조금 바꾼아이디어.
- Gradient 에 velocity 를 가중하 여 계산하는것이 기존방법
- Nesterov는 현재지점에서 velocity를 먼저 계산한 후에 그 지점에서 gradient를 계산 후 원 점으로 돌아가 합함. Velocity값 을 그래디언트 값으로 보정함.
- Velocity는 초기값을 항상 0으로 함.(Not hyperparameter)



단점은 convex 하게 최적화된곳에서는 성능이 좋지만 Neural Network 처럼 non convex한 환경에서는 그다지 성능을 보여주지 못함.

Nesterov Momentum



$$v_{t+1} = \rho v_t - \alpha \nabla f(x_t + \rho v_t)$$
$$x_{t+1} = x_t + v_{t+1}$$

Change of variables $\tilde{x}_t = x_t + \rho v_t$ and rearrange:

$$v_{t+1} = \rho v_t - \alpha \nabla f(\tilde{x}_t)$$

$$\tilde{x}_{t+1} = \tilde{x}_t - \rho v_t + (1+\rho)v_{t+1}$$

$$= \tilde{x}_t + v_{t+1} + \rho(v_{t+1} - v_t)$$

Nesterov는 속도에 에러값을 보정하고 local minima를 지나 global minima(이 경우)로 수렴하게 해주고 overshooting 도 줄여줌.

AdaGrad

- Scalaing개념. Gradient의 제곱을 계속더한값으로 나눠줌.
- step1) $G = G_0^2$, $update \frac{1}{\sqrt{G}}G_0$ step2) $G = G_0^2 + G_1^2$, $update \frac{1}{\sqrt{G}}G_1$... over and over
- Problem is G 값이 너무 커짐. 점점 업데이트 속도가 느려짐.
- → Convex에서는 괜찮지만 non convex의 경우 local minima, saddle point에 안착하게됨.

RMSProp

- AdaGrad의 대안
- 아이디어는 유지하면서

Local minima 등에 안빠지게

속도를 조절!

Decay rate 를 곱해주자!

Almost decay = 0.9 or 0.99

AdaGrad

grad_squared = 0

 $dx = compute_gradient(x)$ grad_squared += dx * dx

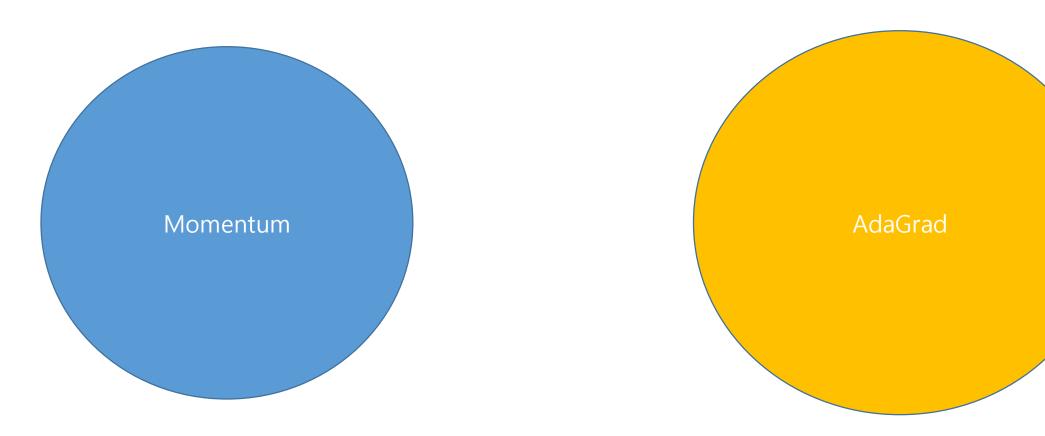
while True:

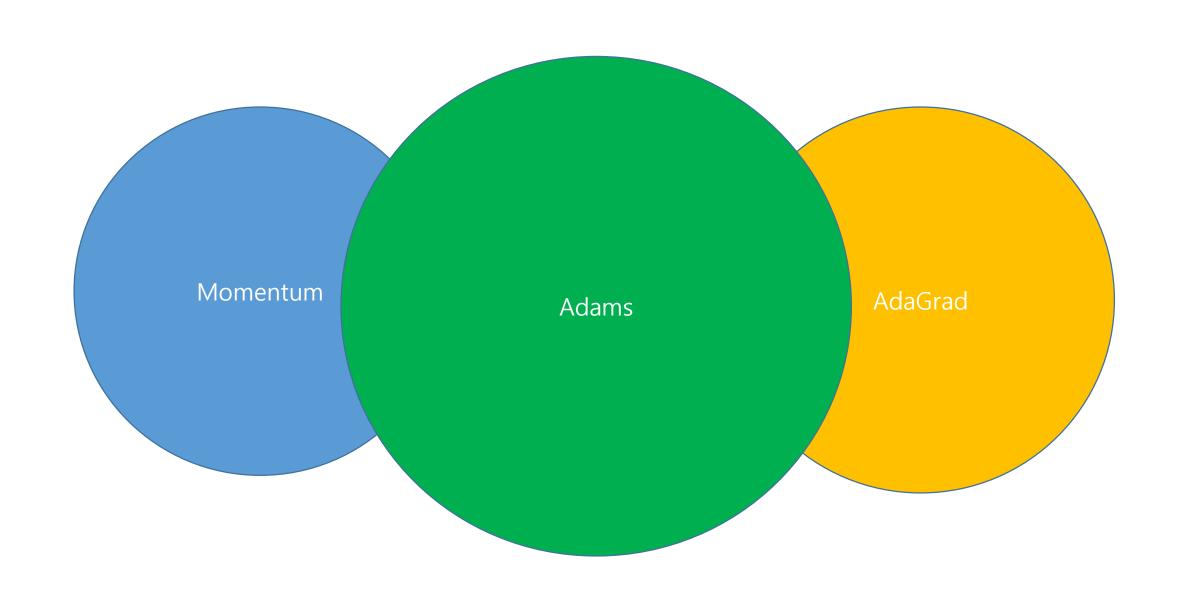
RMSProp

```
grad_squared = 0
while True:
 dx = compute_gradient(x)
 grad_squared = decay_rate * grad_squared + (1 - decay_rate) * dx * dx
  x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)
```

x -= learning_rate * dx / (np.sqrt(grad_squared) + 1e-7)

Optimization....
Two types idea, and still both are good idea





Adams(Almost)

```
first_moment = 0
second_moment = 0
while True:
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx
    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx
    x -= learning_rate * first_moment / (np.sqrt(second_moment) + 1e-7))
Momentum

AdaGrad / RMSProp
```

Sort of like RMSProp with momentum

작은 문제가 있음. second momen가 처음에 너무 작은값이라 step 이 너무 큼.

조정이 필요함

Adam(full form)

```
first_moment = 0
second_moment = 0
for t in range(num_iterations):
    dx = compute_gradient(x)
    first_moment = beta1 * first_moment + (1 - beta1) * dx

    second_moment = beta2 * second_moment + (1 - beta2) * dx * dx

first_unbias = first_moment / (1 - beta1 ** t)
    second_unbias = second_moment / (1 - beta2 ** t)

x -= learning_rate * first_unbias / (np.sqrt(second_unbias) + 1e-7))

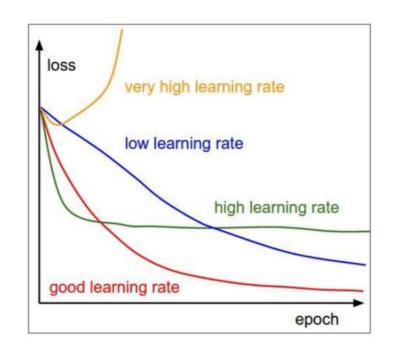
AdaGrad / RMSProp
```

Bias correction for the fact that first and second moment estimates start at zero

Adam with beta1 = 0.9, beta2 = 0.999, and learning_rate = 1e-3 or 5e-4 is a great starting point for many models!

Learning rate..

- SGD, AdaGrad, Momentum, Adam,, all have learning rate term.
- Learning rate is hyperparameter. 하나로 고정하지 않음.



=> Learning rate decay over time!

step decay:

e.g. decay learning rate by half every few epochs.

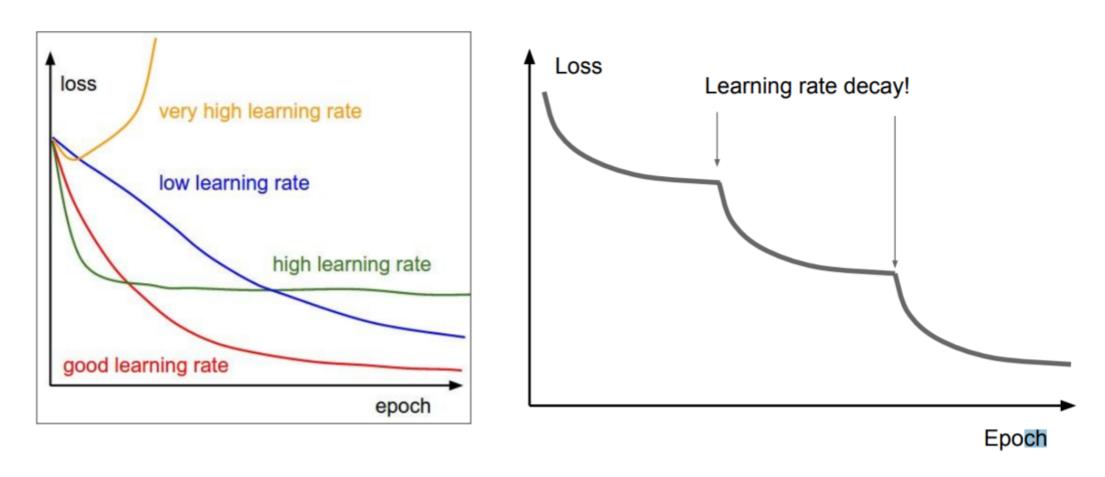
exponential decay:

$$\alpha = \alpha_0 e^{-kt}$$

1/t decay:

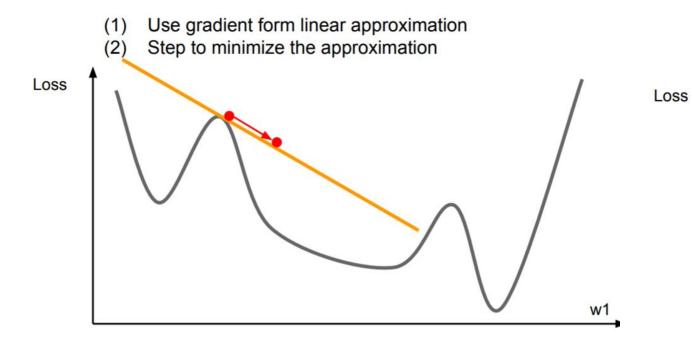
$$lpha=lpha_0/(1+kt)$$

From ResNet



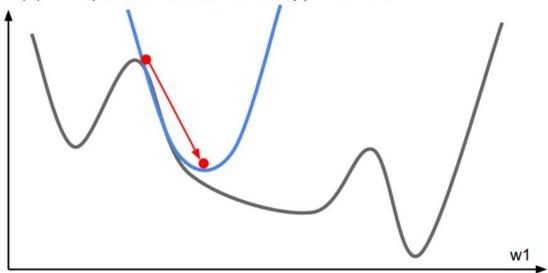
Decay가있다고 하더라도 초기화는 중요하다. 처음에 잘선택하자.

First-Order and Second-Order optimizer



최종적으로는 learning rate 가 필요하지만 지역적으로는 필요없음. 항상 로컬 미니멈을 찾음. 하지만 hessian matrix는 계산 비용이 너무큼(inverse $O(N^3)$)

- (1) Use gradient and Hessian to form quadratic approximation
- (2) Step to the **minima** of the approximation



second-order Taylor expansion:

$$J(\boldsymbol{\theta}) \approx J(\boldsymbol{\theta}_0) + (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0) + \frac{1}{2} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)^{\top} \boldsymbol{H} (\boldsymbol{\theta} - \boldsymbol{\theta}_0)$$

$$\boldsymbol{\theta}^* = \boldsymbol{\theta}_0 - \boldsymbol{H}^{-1} \nabla_{\boldsymbol{\theta}} J(\boldsymbol{\theta}_0)$$

L-BFGS

• Hessian 을 근사시켜 계산비용을 조금 줄임.

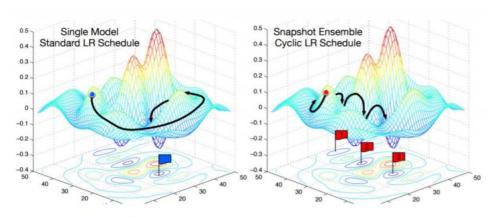
• 그래도 deep neural network에서는 잘 안씀. 조금 변형에서 간 간히 쓰일곳이있음. → 확률적인 모델을 잘 다룰수가없음. Nonconvex에서는 성능이 별로임.

Beyond Training Error

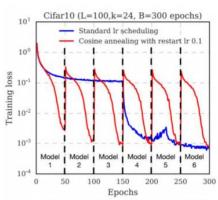
- Training error를 줄이기 위해 노력했음. 하지만 결국 중요한건 한번도 보지않은 데이터에 대한 결과임.(unseen data)
- Train Data와 real data(unseen data)의 갭을 어떻게 줄일까?

- Model Ensembles
 - Train Model independently and ensemble.

Tricks and tips



Loshchilov and Hutter, "SGDR: Stochastic gradient descent with restarts", arXiv 2016 Huang et al, "Snapshot ensembles: train 1, get M for free", ICLR 2017 Figures copyright Yixuan Li and Geoff Pleiss, 2017. Reproduced with permission.



Cyclic learning rate schedules can make this work even better!

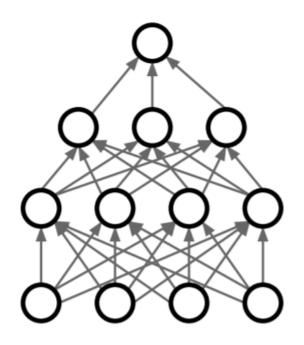
 여러 개 모델을 돌리는 대 신 하나의 모델을 여러 스 냅샷으로 트레이닝.

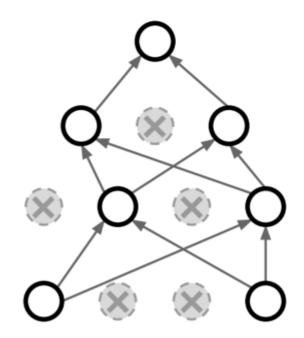
```
while True:
   data_batch = dataset.sample_data_batch()
   loss = network.forward(data_batch)
   dx = network.backward()
   x += - learning_rate * dx
   x_test = 0.995*x_test + 0.005*x # use for test set
```

 실제 파라미터를 사용하지 않고 파라미터의 평균을 계속 갱신해 사용. 조금 성 능 향상. 잘 안씀.

No ensemble, single model

- Multiple model ensemble 은 효율이 떨어짐. 어떻게 성능을 올릴까.? → Regularization!
- 기존에 배운 L2,L1은 NN에서는 별로 좋진않음.
- Drop out 이라는 방법을 사용





Forward pass 에서 임의의 뉴런 을 0으로 만들자.

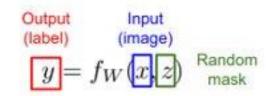
- 1. Overfitting 을 막아줌.
- 2. 매번 다른 뉴런을 drop out 하기 때문에 여러 모델을 앙 상블한 효과.

Test Time에서 랜덤이있는건 별로 좋 지않다. 고양이를 어제는 고양이로 오늘은 개로 인식하면 매우 이상할것.

기댓값을 average out 시켜줘야함.

Dropout: Test time

Dropout makes our output random!



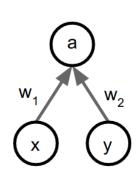
Want to "average out" the randomness at test-time

$$y = f(x) = E_z[f(x,z)] = \int p(z)f(x,z)dz$$

But this integral seems hard ...

Want to approximate the integral

$$y = f(x) = E_z[f(x, z)] = \int p(z)f(x, z)dz$$



Consider a single neuron.

by dropout probability

At test time we have: $E[a] = w_1x + w_2y$ During training we have: $E[a] = \frac{1}{4}(w_1x + w_2y) + \frac{1}{4}(w_1x + 0y)$ $+\frac{1}{4}(0x+0y)+\frac{1}{4}(0x+w_2y)$ At test time, **multiply** $=\frac{1}{2}(w_1x+w_2y)$

```
""" Vanilla Dropout: Not recommended implementation (see notes below) """
                                                                        Dropout Summary
p = 0.5 # probability of keeping a unit active. higher = less dropout
def train_step(X):
 """ X contains the data """
 # forward pass for example 3-layer neural network
 H1 = np.maximum(0, np.dot(W1, X) + b1)
 U1 = np.random.rand(*H1.shape) < p # first dropout mask
 H1 *= U1 # drop!
                                                                           drop in forward pass
 H2 = np.maximum(0, np.dot(W2, H1) + b2)
 U2 = np.random.rand(*H2.shape) < p # second dropout mask
 H2 *= U2 # drop!
 out = np.dot(W3, H2) + b3
 # backward pass: compute gradients... (not shown)
 # perform parameter update... (not shown)
def predict(X):
 # ensembled forward pass
                                                                            scale at test time
 H1 = np.maximum(0, np.dot(W1, X) + b1) * p # NOTE: scale the activations
 H2 = np.maximum(0, np.dot(W2, H1) + b2) * p # NOTE: scale the activations
 out = np.dot(W3, H2) + b3
```

Test time은 성능을 빠르게 하고 싶다.

```
p = 0.5 # probability of keeping a unit active. higher = less dropout
def train step(X):
  # forward pass for example 3-layer neural network
 H1 = np.maximum(0, np.dot(W1, X) + b1)
 U1 = (np.random.rand(*H1.shape) < p) / p # first dropout mask. Notice /p!
  H1 *= U1 # drop!
  H2 = np.maximum(0, np.dot(W2, H1) + b2)
  U2 = (np.random.rand(*H2.shape) < p) / p # second dropout mask. Notice /p!
  H2 *= U2 # drop!
  out = np.dot(W3, H2) + b3
  # backward pass: compute gradients... (not shown)
  # perform parameter update... (not shown)
def predict(X):
  # ensembled forward pass
  H1 = np.maximum(0, np.dot(W1, X) + b1) # no scaling necessary
  H2 = np.maximum(0, np.dot(W2, H1) + b2)
  out = np.dot(W3, H2) + b3
```

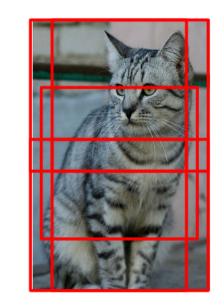
Train 에서 미리 반영. → Drop out을 사용하면 Train시간이 늘어남.

test time is unchanged!

Data Augmentation Random crops and scales

Training: sample random crops / scales ResNet:

- 1. Pick random L in range [256, 480]
- 2. Resize training image, short side = L
- 3. Sample random 224 x 224 patch



Testing: average a fixed set of crops

ResNet:

- 1. Resize image at 5 scales: {224, 256, 384, 480, 640}
- 2. For each size, use 10 224 x 224 crops: 4 corners + center, + flips

Regularization 효과.

Common Pattern

Batch norm이 가장일반적임. 이걸로 regularization이 모자르다면 drop out 추가. 그 이상은 잘없음.

Regularization: A common pattern

Training: Add random noise

Testing: Marginalize over the noise

Examples:

Dropout

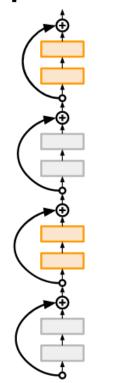
Batch Normalization

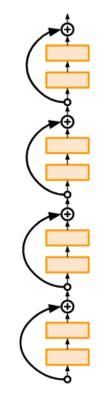
Data Augmentation

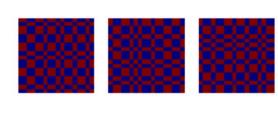
DropConnect

Fractional Max Pooling

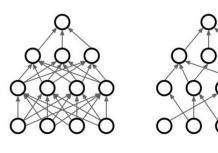
Stochastic Depth







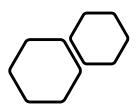
Fractional Max pooling



DropConnection

Huang et al, "Deep Networks with Stochastic Depth", ECCV 2016

Stochastic Depth



Thank you

Q&A