Lecture 12

Common Optimization Algorithms

STAT 479: Deep Learning, Spring 2019 Sebastian Raschka <u>http://stat.wisc.edu/~sraschka/teaching/stat479-ss2019/</u>

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Overview: Additional Tricks for Neural Network Training (Part 2/2)

Part 1 (before Spring break)

- Input Normalization (BatchNorm, InstanceNorm, GroupNorm, LayerNorm)
- Weight Initialization (Xavier, Kaiming He)

Part 2 (this lecture)

- Learning Rate Decay
- Momentum Learning
- Adaptive Learning

Overview: Additional Tricks for Neural Network Training (Part 2/2)

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- Input Normalization (BatchNorm, InstanceNorm, GroupNorm, LayerNorm)
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Part 2 (this lecture)

- Learning Rate Decay
- Momentum Learning
- (Modifications of the 1st order SGD optimization algorithm; 2nd order methods are rarely used in DL)

• Adaptive Learning

Minibatch Learning Recap



 Main advantage: Convergence speed, because it offers to opportunities for parallelism (<u>do you recall what these are?</u>)

- Minibatch learning is a form of stochastic gradient descent
- Each minibatch can be considered a sample drawn from the training set (where the training set is in turn a sample drawn from the population)
- Hence, the gradient is noisier
- A noisy gradient can be
 - ✤ good: chance to escape local minima
 - \blacklozenge bad: can lead to extensive oscillation

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- A noisy gradient can be
 - ✤ good: chance to escape local minima
 - \blacklozenge bad: can lead to extensive oscillation
 - Note that second order methods that take e.g., gradient curvature into account usually don't work so well in practice and are not often used/recommended in DL

Nice Library & Visualization Tool

https://vis.ensmallen.org



Practical Tip for Minibatch Use

- Reasonable minibatch sizes are usually: 32, 64, 128, 256, 512, 1024 (in the last lecture, we discussed why powers of 2 are a common convention)
- Usually, you can choose a batch size that is as large as your GPU memory allows (matrix-multiplication and the size of fully-connected layers are usually the bottleneck)
- Practical tip: usually, it is a good idea to also make the batch size proportional to the number of classes in the dataset



Raschka, S. (2018). Model evaluation, model selection, and algorithm selection in machine learning. *arXiv preprint arXiv:1811.12808*.

Figure 1: Distribution of Iris flower classes upon random subsampling into training and test sets.

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- Batch effects -- minibatches are samples of the training set, hence minibatch loss and gradients are approximations
- Hence, we usually get oscillations
- To dampen oscillations towards the end of the training, we can decay the learning rate

exponentially weighted average or whole-training set loss minibato loss Mannammann i Levelion

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Loss

- Batch effects -- minibatches are samples of the training set, hence minibatch loss and gradients are approximations
- Hence, we usually get oscillations
- To dampen oscillations towards the end of the training, we can decay the learning rate

loss

Danger of learning rate is to decrease the learning rate too early

Practical tip: try to train the model without learning rate decay first, then add it later

You can also use the validation performance (e.g., accuracy) to judge whether Ir decay is useful (as opposed to using the training loss)

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Most common variants for learning rate decay:

1) Exponential Decay:

 $\eta_t := \eta_0 \cdot e^{-k \cdot t}$

where k is the decay rate



Most common variants for learning rate decay:

2) Halving the learning rate:

 $\eta_t := \eta_{t=1}/2$

3) Inverse decay:

$$\eta_t := \frac{\eta_0}{1 + k \cdot t}$$



There are many, many more

E.g., Cyclical Learning Rate

Smith, Leslie N. "Cyclical learning rates for training neural networks." Applications of Computer Vision (WACV), 2017 IEEE Winter Conference on. IEEE, 2017.



Figure 2. Triangular learning rate policy. The blue lines represent learning rate values changing between bounds. The input parameter *stepsize* is the number of iterations in half a cycle.

(which, I found, didn't work well at all in practice, unfortunately -- at least in my case)

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Geoffrey E Hinton



Yann LeCun



Yoshua Bengio



FATHERS OF THE DEEP LEARNING REVOLUTION RECEIVE ACM A.M. TURING AWARD

Bengio, Hinton, and LeCun Ushered in Major Breakthroughs in Artificial Intelligence

ACM named Yoshua Bengio, Geoffrey Hinton, and Yann LeCun recipients of the 2018 ACM A.M. Turing Award for conceptual and engineering breakthroughs that have made deep neural networks a critical component of computing. Bengio is Professor at the University of Montreal and Scientific Director at Mila, Quebec's Artificial Intelligence Institute; Hinton is VP and Engineering Fellow of Google, Chief Scientific Adviser of The Vector Institute, and University Professor Emeritus at the University of Toronto; and LeCun is Professor at New York University and VP and Chief AI Scientist at Facebook.

Working independently and together, Hinton, LeCun and Bengio developed conceptual foundations for the field, identified surprising phenomena through experiments, and contributed engineering advances that demonstrated the practical advantages of deep neural networks. In recent years, deep learning methods have been responsible for astonishing breakthroughs in computer vision, speech recognition, natural language processing, and robotics among other applications.

While the use of artificial neural networks as a tool to help computers recognize patterns and simulate human intelligence had been introduced in the 1980s, by the early 2000s, LeCun, Hinton and Bengio were among a small group who remained committed to this approach. Though their efforts to rekindle the AI community's interest in neural networks were initially met with skepticism, their ideas recently resulted in major technological advances, and their methodology is now the dominant paradigm in the field.

Some Live News

The Turing Award is generally recognized as the highest distinction in computer science and the "Nobel Prize of computing". ... Since 2014, the award has been accompanied by a prize of US \$1 million

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<u>Option 1.</u> Just call your own function at the end of each epoch:

```
def adjust_learning_rate(optimizer, epoch, initial_lr, decay_rate):
    """Exponential decay every 10 epochs"""
    if not epoch % 10:
        lr = initial_lr * torch.exp(-decay_rate*epoch)
        for param_group in optimizer.param_groups:
            param group['lr'] = lr
```

<u>Option 2.</u> Use one of the built-in tools in PyTorch: (many more available) (Here, the most generic version.)

CLASS torch.optim.lr_scheduler.LambdaLR(*optimizer*, *lr_lambda*, *last_epoch=-1*)

Sets the learning rate of each parameter group to the initial Ir times a given function. When last_epoch=-1, sets initial Ir as Ir.

Parameters:

- optimizer (Optimizer) Wrapped optimizer.
 - Ir_lambda (function or list) A function which computes a multiplicative factor given an integer parameter epoch, or a list of such functions, one for each group in optimizer.param_groups.
 - last_epoch (int) The index of last epoch. Default: -1.

Example

```
>>> # Assuming optimizer has two groups.
>>> lambda1 = lambda epoch: epoch // 30
>>> lambda2 = lambda epoch: 0.95 ** epoch
>>> scheduler = LambdaLR(optimizer, lr_lambda=[lambda1, lambda2])
>>> for epoch in range(100):
>>> scheduler.step()
>>> train(...)
>>> validate(...) Source: https://pytorch.org/docs/stable/optim.html
```

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[SOURCE]

```
model = model.to(DEVICE)
```

```
optimizer = torch.optim.SGD(model.parameters(), lr=0.1)
```

https://github.com/rasbt/stat479-deep-learning-ss19/tree/master/L12_optim/lr_scheduler_and_saving_models.ipynb

Example, part 1/2

```
for epoch in range(5):
   model.train()
                                                                                 Example, part 2/2
   for batch idx, (features, targets) in enumerate(train loader):
       features = features.view(-1, 28*28).to(DEVICE)
       targets = targets.to(DEVICE)
       ### FORWARD AND BACK PROP
       logits, probas = model(features)
       #cost = F.nll_loss(torch.log(probas), targets)
       cost = F.cross entropy(logits, targets)
       optimizer.zero grad()
       cost.backward()
       minibatch cost.append(cost)
       ### UPDATE MODEL PARAMETERS
       optimizer.step()
       ### LOGGING
       if not batch idx % 50:
           print ('Epoch: %03d/%03d | Batch %03d/%03d | Cost: %.4f'
                  %(epoch+1, NUM EPOCHS, batch idx,
                   len(train loader), cost))
   ### Update Learning Rate
   scheduler.step() # don't have to do it every epoch!
```

model.eval()

https://github.com/rasbt/stat479-deep-learning-ss19/tree/master/L12_optim/lr_scheduler_and_saving_models.ipynb

Saving Models in PyTorch

Save Model

```
model.to(torch.device('cpu'))
torch.save(model.state_dict(), './my_model_2epochs.pt')
torch.save(optimizer.state_dict(), './my_optimizer_2epochs.pt')
torch.save(scheduler.state_dict(), './my_scheduler_2epochs.pt')
```

Load Model

Learning rate schedulers have the advantage that we can also simply save their state for reuse (e.g., saving and continuing training later)

<u>https://github.com/rasbt/stat479-deep-learning-ss19/tree/master/L12_optim/lr_scheduler_and_saving_models.ipynb</u>

Weight Initialization Experiments (Last-lecture-follow-up)

https://github.com/rasbt/stat479-deep-learning-ss19/tree/master/L13_intro-cnn/code/cnn-with-diff-init_

Uniform: Test accuracy 97.63%

def weights_init(m):

if isinstance(m, nn.Linear) or isinstance(m, nn.Conv2d):
 torch.nn.init.uniform_(m.weight.detach(), -0.1, 0.1)
 torch.zero_(m.bias.detach())

model.apply(weights_init)

Default: Test accuracy 97.77%

Xavier Normal: Test accuracy 97.69%

def weights_init(m):

if isinstance(m, nn.Linear) or isinstance(m, nn.Conv2d):
 torch.nn.init.xavier_normal_(m.weight)
 torch.zero_(m.bias.detach())

model.apply(weights_init)

Normal: Test accuracy 97.76%

def weights_init(m):

if isinstance(m, nn.Linear) or isinstance(m, nn.Conv2d):
 torch.nn.init.normal_(m.weight.detach(), mean=0, std=0.1)
 torch.zero_(m.bias.detach())

model.apply(weights_init)

Xavier Uniform: Test accuracy 97.36%

def weights_init(m):

if isinstance(m, nn.Linear) or isinstance(m, nn.Conv2d):
 torch.nn.init.xavier_uniform_(m.weight)
 torch.zero_(m.bias.detach())

model.apply(weights_init)

He Normal: Test accuracy 97.67%

def weights_init(m):

if isinstance(m, nn.Linear) or isinstance(m, nn.Conv2d):
 torch.nn.init.kaiming_normal_(m.weight)
 torch.zero_(m.bias.detach())

model.apply(weights_init)

He Uniform: Test accuracy 97.54%

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```
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```

Momentum

From Wikipedia, the free encyclopedia

This article is about linear momentum. It is not to be confused with angular momentum. This article is about momentum in physics. For other uses, see Momentum (disambigua

In Newtonian mechanics, linear momentum, translational momentum, or simply momentum (pl. momenta) is the product of the mass and velocity of an object. It is a vector quantity, possessing a magnitude and a direction in three-dimensional space. If m is an object's mass and v is the velocity (also a vector), then the momentum is

Source: https://en.wikipedia.org/wiki/Momentum

- Momentum is a jargon term in DL and is probably a misnomer in this context
- Concept: In momentum learning, we try to accelerate convergence by dampening oscillations using "velocity" (the speed of the "movement" from previous updates)

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Without momentum

With momentum







With momentum

Key take-away:

Not only move in the (opposite) direction of the gradient, but also move in the "averaged" direction of the last few updates

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Not only move in the (opposite) direction of the gradient, but also move in the "averaged" direction of the last few updates

Helps with dampening oscillations, but also helps with escaping local minima traps





Weight update using the velocity vector:

$$w_{i,j}(t+1) := w_{i,j}(t) - \Delta w_{i,j}(t)$$

Qian, N. (1999). On the momentum term in gradient descent learning algorithms. Neural Networks : The Official Journal of the International Neural Network Society, 12(1), 145–151. http://doi.org/10.1016/S0893-6080(98)00116-6





We often think of Momentum as a means of dampening oscillations and speeding up the iterations, leading to faster convergence. But it has other interesting behavior. It allows a larger range of step-sizes to be used, and creates its own oscillations. What is going on?



Step-size a = 0.02			Momentum $\beta = 0.99$		
1 0	0.003	0.006	0.00	0.500	0.990

and speeding up the iterations, leading to faster convergence. But it has other interesting behavior. It allows a larger range of step-sizes to be used, and creates its own oscillations. What is going on?

Source: https://distill.pub/2017/momentum/

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[SOURCE]

Implements stochastic gradient descent (optionally with momentum).

Nesterov momentum is based on the formula from On the importance of initialization and momentum in deep learning.

Parameters:

- params (iterable) iterable of parameters to optimize or dicts defining parameter groups
 - Ir (float) learning rate
 - momentum (float, optional) momentum factor (default: 0)
 - weight_decay (float, optional) weight decay (L2 penalty) (default: 0)
 - dampening (float, optional) dampening for momentum (default: 0)
 - nesterov (bool, optional) enables Nesterov momentum (default: False)

Example

Source: <u>https://pytorch.org/docs/stable/optim.html</u>

[SOURCE]

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Example

Note that the optional "dampening" term is used as follows:

```
v = momentum * v + (1-dampening) * gradientW
W = W - lr * v
```

Also note that in PyTorch, the learning rate is also applied to the momentum terms, instead of the original definition, which would be

```
v = momentum * v + (1-dampening) * lr * gradientW
W = W - v
```

A Better Momentum Method: Nesterov Accelerated Gradient

Similar to momentum learning, but performs a correction after the update (based on where the loss, w.r.t. the weight parameters, is approx. <u>going to be after the update</u>)

$$\begin{array}{l} \underline{\text{Before:}}\\ \Delta \mathbf{w}_t := \alpha \cdot \Delta \mathbf{w}_{t-1} + \eta \cdot \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w}_t)\\ \mathbf{w}_{t+1} := \mathbf{w}_t - \Delta \mathbf{w}_t \end{array}$$

Nesterov:

$$\Delta \mathbf{w}_t := \alpha \cdot \Delta \mathbf{w}_{t-1} + \eta \cdot \nabla_{\mathbf{w}} \mathcal{L}(\mathbf{w}_t - \alpha \cdot \Delta \mathbf{w}_{t-1})$$
$$\mathbf{w}_{t+1} := \mathbf{w}_t - \Delta \mathbf{w}_t$$

Nesterov, Y. (1983). A method for unconstrained convex minimization problem with the rate of convergence $o(1/k^2)$. Doklady ANSSSR (translated as Soviet.Math.Docl.), vol. 269, pp. 543– 547.

Sutskever, I., Martens, J., Dahl, G. E., & Hinton, G. E. (2013). On the importance of initialization and momentum in deep learning. *ICML (3)*, *28*(1139-1147), 5.

A Better Momentum Method: Nesterov Accelerated Gradient



Figure 1. (Top) Classical Momentum (Bottom) Nesterov Accelerated Gradient

Sutskever, I., Martens, J., Dahl, G. E., & Hinton, G. E. (2013). On the importance of initialization and momentum in deep learning. *ICML (3)*, *28*(1139-1147), 5.

A Better Momentum Method: Nesterov Accelerated Gradient



Sutskever, I., Martens, J., Dahl, G. E., & Hinton, G. E. (2013). On the importance of initialization and momentum in deep learning. *ICML (3)*, *28*(1139-1147), 5.

Adaptive Learning Rates

There are many different flavors of adapting the learning rate (bit out of scope for this course to review them all)

<u>Key take-aways:</u>

- decrease learning if the gradient changes its direction
- increase learning if the gradient stays consistent

Adaptive Learning Rates

Key take-aways:

- decrease learning if the gradient changes its direction
- increase learning if the gradient stays consistent

Step 1: Define a local gain (g) for each weight (initialized with g=1)

$$\Delta w_{i,j} := \eta \cdot g_{i,j} \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}$$

Adaptive Learning Rates

<u>Step 1</u>: Define a local gain (g) for each weight (initialized with g=1)

$$\Delta w_{i,j} := \eta \cdot g_{i,j} \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}$$

<u>Step 2:</u>

If gradient is consistent $g_{i,j}(t) := g_{i,j}(t-1) + \beta$ else

$$g_{i,j}(t) := g_{i,j}(t-1) \cdot (1-\beta)$$

Note that

multiplying by a factor has a larger impact if gains are large, compared to adding a term

(dampening effect if updates oscillate in the wrong direction)

Adaptive Learning Rate via RMSProp

- Unpublished algorithm by Geoff Hinton (but very popular) based on Rprop [1]
- Very similar to another concept called AdaDelta
- Concept: divide learning rate by exponentially decreasing moving average of the squared gradients
- This takes into account that gradients can vary widely in magnitude
- Here, RMS stands for "Root Mean Squared"
- Also, damps oscillations like momentum (but in practice, works a bit better)

[1] Igel, Christian, and Michael Hüsken. "Improving the Rprop learning algorithm." *Proceedings of the Second International ICSC Symposium on Neural Computation (NC 2000)*. Vol. 2000. ICSC Academic Press, 2000.

Adaptive Learning Rate via RMSProp

$$MeanSquare(w_{i,j}, t) := \beta \cdot MeanSquare(w_{i,j}, t-1) + (1-\beta) \left(\frac{\partial \mathcal{L}}{w_{i,j}(t)}\right)^{2}$$

moving average of the squared gradient for each weight

$$w_{i,j}(t) := w_{i,j}(t) - \eta \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}(t)} / \left(\sqrt{MeanSquare\left(w_{i,j}, t\right)} + \epsilon \right)$$

where beta is typically between 0.9 and 0.999

small epsilon term to avoid division by zero

Adaptive Learning Rate via ADAM

- ADAM (Adaptive Moment Estimation) is probably the most widely used optimization algorithm in DL as of today
- It is a combination of the momentum method and RMSProp

$$\begin{array}{c} \underbrace{\text{Momentum-like term:}}_{\Delta w_{i,j}(t) := \alpha \cdot \Delta w_{i,j}(t-1) + \eta \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}(t)}_{\Delta w_{i,j}(t) := \alpha \cdot \Delta w_{i,j}(t-1) + \eta \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}(t)} \text{ original momentum term} \\ m_t \\ m_t := \alpha \cdot m_{t-1} + (1-\alpha) \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}(t) \\ \underbrace{\text{RMSProp term:}}_{MeanSquare}(w_{i,j}, t) := \beta \cdot MeanSquare(w_{i,j}, t-1) + (1-\beta) \left(\frac{\partial \mathcal{L}}{\partial w_{i,j}(t)}\right)^2 \end{array}$$

Kingma, D. P., & Ba, J. (2014). Adam: A method for stochastic optimization. arXiv preprint arXiv:1412.6980.

Adaptive Learning Rate via ADAM

Momentum-like term:

$$\overline{\Delta w_{i,j}(t)} := \alpha \cdot \Delta w_{i,j}(t-1) + \eta \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}(t) \quad \text{original momentum term} \\
m_t := \alpha \cdot m_{t-1} + (1-\alpha) \cdot \frac{\partial \mathcal{L}}{\partial w_{i,j}}(t)$$

$$\frac{RMSProp \text{ term:}}{r := \beta \cdot MeanSquare(w_{i,j}, t-1) + (1-\beta) \left(\frac{\partial \mathcal{L}}{\partial w_{i,j}(t)}\right)^2$$

ADAM update:

$$w_{i,j} := w_{i,j} - \eta \frac{m_t}{\sqrt{r} + \epsilon}$$

Kingma, D. P., & Ba, J. (2014). Adam: A method for stochastic optimization. arXiv preprint arXiv:1412.6980.

Adaptive Learning Rate via ADAM

Algorithm 1: Adam, our proposed algorithm for stochastic optimization. See section 2 for details, and for a slightly more efficient (but less clear) order of computation. g_t^2 indicates the elementwise square $g_t \odot g_t$. Good default settings for the tested machine learning problems are $\alpha = 0.001$, $\beta_1 = 0.9$, $\beta_2 = 0.999$ and $\epsilon = 10^{-8}$. All operations on vectors are element-wise. With β_1^t and β_2^t we denote β_1 and β_2 to the power t.

Require: α : Stepsize **Require:** $\beta_1, \beta_2 \in [0, 1)$: Exponential decay rates for the moment estimates **Require:** $f(\theta)$: Stochastic objective function with parameters θ **Require:** θ_0 : Initial parameter vector $m_0 \leftarrow 0$ (Initialize 1st moment vector) $v_0 \leftarrow 0$ (Initialize 2nd moment vector) $t \leftarrow 0$ (Initialize timestep) while θ_t not converged **do** $t \leftarrow t + 1$ $g_t \leftarrow \nabla_{\theta} f_t(\theta_{t-1})$ (Get gradients w.r.t. stochastic objective at timestep t) $m_t \leftarrow \beta_1 \cdot m_{t-1} + (1 - \beta_1) \cdot g_t$ (Update biased first moment estimate) $v_t \leftarrow \beta_2 \cdot v_{t-1} + (1 - \beta_2) \cdot g_t^2$ (Update biased second raw moment estimate) $\widehat{m}_t \leftarrow m_t/(1 - \beta_1^t)$ (Compute bias-corrected first moment estimate) $\hat{v}_t \leftarrow v_t/(1-\beta_2^t)$ (Compute bias-corrected second raw moment estimate) $\theta_t \leftarrow \theta_{t-1} - \alpha \cdot \hat{m}_t / (\sqrt{\hat{v}_t} + \epsilon)$ (Update parameters) Also add a bias correction term end while for better conditioning in earlier iterations **return** θ_t (Resulting parameters)

Kingma, D. P., & Ba, J. (2014). Adam: A method for stochastic optimization. *arXiv preprint arXiv:1412.6980*.



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https://bl.ocks.org/EmilienDupont/aaf429be5705b219aaaf8d691e27ca87

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Using Different Optimizers in PyTorch

Usage is the as for vanilla SGD, which we used before,

you can find an overview at: <u>https://pytorch.org/docs/stable/optim.html</u>

optimizer = torch.optim.SGD(model.parameters(), lr=0.01, momentum=0.9)
optimizer = torch.optim.Adam(model.parameters(), lr=0.0001)

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Usage is the as for vanilla SGD, which we used before,

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optimizer = torch.optim.SGD(model.parameters(), lr=0.01, momentum=0.9)
optimizer = torch.optim.Adam(model.parameters(), lr=0.0001)

Remember to save the optimizer state if you are using, e.g., Momentum or ADAM, and want to continue training later (see earlier slides on saving states of the learning rate schedulers).

Training Loss vs Generalization Error

Improving Generalization Performance by Switching from Adam to SGD

Nitish Shirish Keskar, Richard Socher

(Submitted on 20 Dec 2017)

Despite superior training outcomes, adaptive optimization methods such as Adam, Adagrad or RMSprop have been found to generalize poorly compared to Stochastic gradient descent (SGD). These methods tend to perform well in the initial portion of training but are outperformed by SGD at later stages of training. We investigate a hybrid strategy that begins training with an adaptive method and switches to SGD when appropriate. Concretely, we propose SWATS, a simple strategy which switches from Adam to SGD when a triggering condition is satisfied. The condition we propose relates to the projection of Adam steps on the gradient subspace. By design, the monitoring process for this condition adds very little overhead and does not increase the number of hyperparameters in the optimizer. We report experiments on several standard benchmarks such as: ResNet, SENet, DenseNet and PyramidNet for the CIFAR-10 and CIFAR-100 data sets, ResNet on the tiny-ImageNet data set and language modeling with recurrent networks on the PTB and WT2 data sets. The results show that our strategy is capable of closing the generalization gap between SGD and Adam on a majority of the tasks.

Training Loss vs Generalization Error



Kingma, D. P., & Ba, J. (2014). Adam: A method for stochastic optimization. *arXiv preprint arXiv: 1412.6980*.

Keskar, N. S., & Socher, R. (2017). Improving generalization performance by switching from adam to sgd. *arXiv preprint arXiv:1712.07628*.

Reading Assignment

"An overview of gradient descent optimization algorithms" by Sebastian Ruder: <u>http://ruder.io/optimizing-gradient-descent/index.html</u>