Stories from modern Machine Learning from the optimization perspective

Daniil Merkulov

Optimization methods. MIPT



Neural network is a function, that takes an input x and current set of weights (parameters) w and predicts some vector as an output. Note, that a variety of feed-forward neural networks could be represented as a series of linear transformations, followed by some nonlinear function (say, ReLU (x) or sigmoid):



Neural network is a function, that takes an input x and current set of weights (parameters) w and predicts some vector as an output. Note, that a variety of feed-forward neural networks could be represented as a series of linear transformations, followed by some nonlinear function (say, ReLU (x) or sigmoid):

$$\mathcal{N}\mathcal{N}(\mathbf{w},x) = \sigma_L \circ w_L \circ \ldots \circ \sigma_1 \circ w_1 \circ x \qquad \mathbf{w} = (W_1, b_1, \ldots W_L, b_L),$$



Neural network is a function, that takes an input x and current set of weights (parameters) w and predicts some vector as an output. Note, that a variety of feed-forward neural networks could be represented as a series of linear transformations, followed by some nonlinear function (say, ReLU (x) or sigmoid):

$$\mathcal{N}\mathcal{N}(\mathbf{w},x) = \sigma_L \circ w_L \circ \ldots \circ \sigma_1 \circ w_1 \circ x \qquad \mathbf{w} = (W_1, b_1, \ldots W_L, b_L),$$

where L is the number of layers, σ_i - non-linear activation function, $w_i = W_i x + b_i$ - linear layer.



Neural network is a function, that takes an input x and current set of weights (parameters) w and predicts some vector as an output. Note, that a variety of feed-forward neural networks could be represented as a series of linear transformations, followed by some nonlinear function (say, ReLU (x) or sigmoid):

$$\mathcal{N}\mathcal{N}(\mathbf{w},x) = \sigma_L \circ w_L \circ \ldots \circ \sigma_1 \circ w_1 \circ x \qquad \mathbf{w} = (W_1, b_1, \ldots W_L, b_L),$$

where L is the number of layers, σ_i - non-linear activation function, $w_i = W_i x + b_i$ - linear layer.

Typically, we aim to find \mathbf{w} in order to solve some problem (let say to be $\mathcal{NN}(\mathbf{w}, x_i) \sim y_i$ for some training data x_i, y_i). In order to do it, we solve the optimization problem:

Neural network is a function, that takes an input x and current set of weights (parameters) w and predicts some vector as an output. Note, that a variety of feed-forward neural networks could be represented as a series of linear transformations, followed by some nonlinear function (say, ReLU (x) or sigmoid):

$$\mathcal{N}\mathcal{N}(\mathbf{w},x) = \sigma_L \circ w_L \circ \ldots \circ \sigma_1 \circ w_1 \circ x \qquad \mathbf{w} = (W_1,b_1,\ldots W_L,b_L),$$

where L is the number of layers, σ_i - non-linear activation function, $w_i = W_i x + b_i$ - linear layer.

Typically, we aim to find \mathbf{w} in order to solve some problem (let say to be $\mathcal{NN}(\mathbf{w}, x_i) \sim y_i$ for some training data x_i, y_i). In order to do it, we solve the optimization problem:

$$L(\mathbf{w}, X, y) \to \min_{\mathbf{w}} \qquad \frac{1}{N} \sum_{i=1}^{N} l(\mathbf{w}, x_i, y_i) \to \min_{\mathbf{w}}$$



Loss functions

In the context of training neural networks, the loss function, denoted by $l(\mathbf{w}, x_i, y_i)$, measures the discrepancy between the predicted output $\mathcal{NN}(\mathbf{w}, x_i)$ and the true output y_i . The choice of the loss function can significantly influence the training process. Common loss functions include:

Mean Squared Error (MSE)

Used primarily for regression tasks. It computes the square of the difference between predicted and true values, averaged over all samples.

$$\mathsf{MSE}(\mathbf{w}, X, y) = \frac{1}{N} \sum_{i=1}^{N} (\mathcal{NN}(\mathbf{w}, x_i) - y_i)^2$$

Cross-Entropy Loss

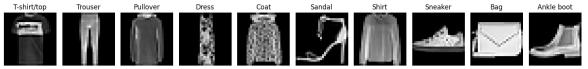
Typically used for classification tasks. It measures the dissimilarity between the true label distribution and the predictions, providing a probabilistic interpretation of classification.

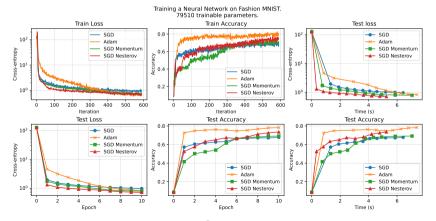
$$\mathsf{Cross-Entropy}(\mathbf{w}, X, y) = -\frac{1}{N} \sum_{i=1}^{N} \sum_{c=1}^{C} y_{i,c} \log(\mathcal{NN}(\mathbf{w}, x_i)_c)$$

where $y_{i,c}$ is a binary indicator (0 or 1) if class label c is the correct classification for observation i, and C is the number of classes.



Simple example: Fashion MNIST classification problem





 $f \rightarrow \min_{x,y,z}$ General introduction

Figure 1: **P**Open in colab

♥ 0 0 4

Visualizing loss surface of neural network via line projection

We denote the initial point as w_0 , representing the weights of the neural network at initialization. The weights after training are denoted as \hat{w} .

Initially, we generate a random Gaussian direction $w_1 \in \mathbb{R}^p$, which inherits the magnitude of the original neural network weights for each parameter group. Subsequently, we sample the training and testing loss surfaces at points along the direction w_1 , situated close to either w_0 or \hat{w} .

Mathematically, this involves evaluating:

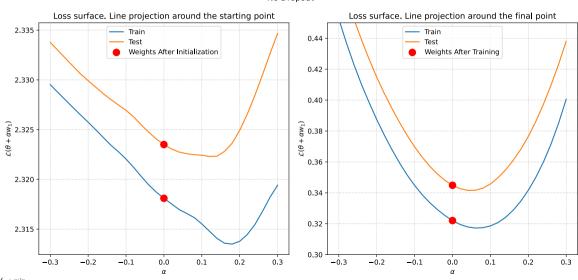
 $L(\alpha) = L(w_0 + \alpha w_1)$, where $\alpha \in [-b, b]$.

Here, α plays the role of a coordinate along the w_1 direction, and b stands for the bounds of interpolation. Visualizing $L(\alpha)$ enables us to project the *p*-dimensional surface onto a one-dimensional axis.

It is important to note that the characteristics of the resulting graph heavily rely on the chosen projection direction. It's not feasible to maintain the entirety of the information when transforming a space with 100,000 dimensions into a one-dimensional line through projection. However, certain properties can still be established. For instance, if $L(\alpha) \mid_{\alpha=0}$ is decreasing, this indicates that the point lies on a slope. Additionally, if the projection is non-convex, it implies that the original surface was not convex.



Visualizing loss surface of neural network

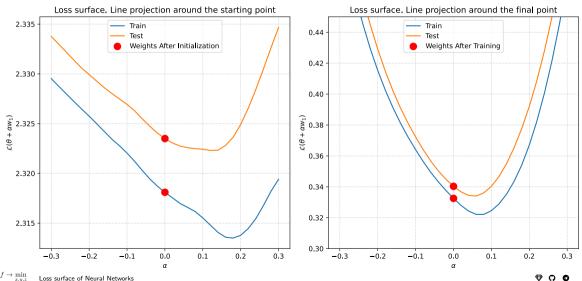


No Dropout

© 0 0 6

 $f \to \min_{x,y,z}$

Visualizing loss surface of neural network



Dropout 0.2

Loss surface of Neural Networks

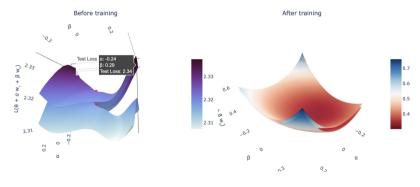
⊕ ∩ ∅ 6

Plane projection

We can explore this idea further and draw the projection of the loss surface to the plane, which is defined by 2 random vectors. Note, that with 2 random gaussian vectors in the huge dimensional space are almost certainly orthogonal. So, as previously, we generate random normalized gaussian vectors $w_1, w_2 \in \mathbb{R}^p$ and evaluate the loss function

$$L(\alpha, \beta) = L(w_0 + \alpha w_1 + \beta w_2)$$
, where $\alpha, \beta \in [-b, b]^2$.

No Dropout. Plane projection of loss surface.



" Train Loss 📲 Test Loss 🔹 Weights before training 👫 Train Loss 👫 Test Loss 🔹 Weights after training



Figure 4: 🏶Open in colab

Can plane projections be useful? ¹

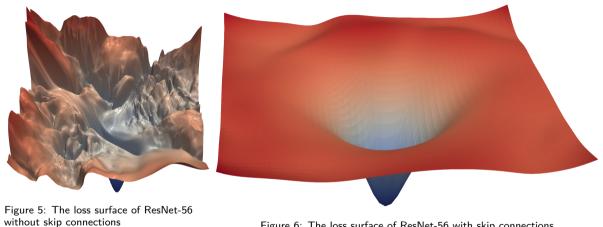


Figure 6: The loss surface of ResNet-56 with skip connections

¹Visualizing the Loss Landscape of Neural Nets, Hao Li, Zheng Xu, Gavin Taylor, Christoph Studer, Tom Goldstein

Can plane projections be useful, really?²

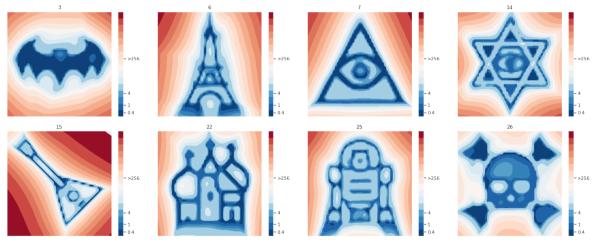


Figure 7: Examples of a loss landscape of a typical CNN model on FashionMNIST and CIFAR10 datasets found with MPO. Loss values are color-coded according to a logarithmic scale

²Loss Landscape Sightseeing with Multi-Point Optimization, Ivan Skorokhodov, Mikhail Burtsev

Impact of initialization ³

- Properly initializing a NN important. NN loss is highly nonconvex; optimizing it to attain a "good" solution hard, requires careful tuning.
- Don't initialize all weights to be the same why?

Impact of initialization ³

- Properly initializing a NN important. NN loss is highly nonconvex; optimizing it to attain a "good" solution hard, requires careful tuning.
- Don't initialize all weights to be the same why?
- Random: Initialize randomly, e.g., via the Gaussian $N(0, \sigma^2)$, where std σ depends on the number of neurons in a given layer. Symmetry breaking.

Impact of initialization³

- Properly initializing a NN important. NN loss is highly nonconvex; optimizing it to attain a "good" solution hard, requires careful tuning.
- Don't initialize all weights to be the same why?
- Random: Initialize randomly, e.g., via the Gaussian $N(0, \sigma^2)$, where std σ depends on the number of neurons in a given layer. Symmetry breaking.
- One can find more useful advices here

³On the importance of initialization and momentum in deep learning Ilya Sutskever, James Martens, George Dahl, Geoffrey Hinton $f \to \min_{x,y,z}$

Impact of initialization ⁴

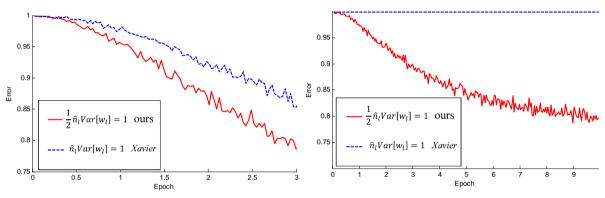


Figure 8: 22-layer ReLU net: good init converges faster

Figure 9: 30-layer ReLU net: good init is able to converge

⁴Delving Deep into Rectifiers: Surpassing Human-Level Performance on ImageNet Classification, Kaiming He, Xiangyu Zhang, Shaoqing Ren, Jian Sun

Grokking ⁵

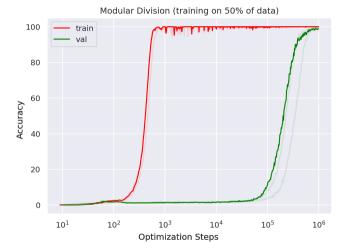
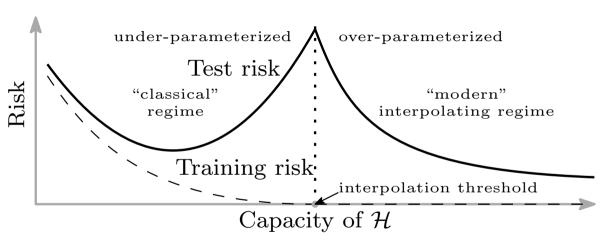


Figure 10: Training transformer with 2 layers, width 128, and 4 attention heads, with a total of about $4 \cdot 10^5$ non-embedding parameters. Reproduction of experiments (~ half an hour) is available here

⁵Grokking: Generalization Beyond Overfitting on Small Algorithmic Datasets, Alethea Power, Yuri Burda, Harri Edwards, Igor Babuschkin, $V_{f \to min}^{f \to min}$ Loss surface of Neural Networks $\heartsuit \circ$ \bullet 12

Double Descent ⁶



13

⁶Reconciling modern machine learning practice and the bias-variance trade-off, Mikhail Belkin, Daniel Hsu, Siyuan Ma, Soumik Mandal $f \rightarrow \min_{x, x, y, z}$ Loss surface of Neural Networks

Exponential learning rate

• Exponential Learning Rate Schedules for Deep Learning

• Multiplication of a chain of matrices in backprop



- Multiplication of a chain of matrices in backprop
- If several of these matrices are "small" (i.e., norms < 1), when we multiply them, the gradient will decrease exponentially fast and tend to vanish (hurting learning in lower layers much more)



- Multiplication of a chain of matrices in backprop
- If several of these matrices are "small" (i.e., norms < 1), when we multiply them, the gradient will decrease exponentially fast and tend to vanish (hurting learning in lower layers much more)
- Conversely, if several matrices have large norm, the gradient will tend to explode. In both cases, the gradients are unstable.

- Multiplication of a chain of matrices in backprop
- If several of these matrices are "small" (i.e., norms < 1), when we multiply them, the gradient will decrease exponentially fast and tend to vanish (hurting learning in lower layers much more)
- Conversely, if several matrices have large norm, the gradient will tend to explode. In both cases, the gradients are unstable.
- Coping with unstable gradients poses several challenges, and must be dealt with to achieve good results.

Feedforward Architecture

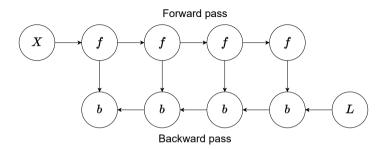


Figure 11: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The activations marked with an f. The gradient of the loss with respect to the activations and parameters marked with b.



Feedforward Architecture

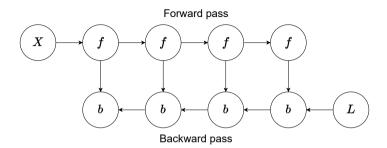


Figure 11: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The activations marked with an f. The gradient of the loss with respect to the activations and parameters marked with b.





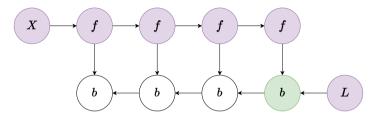


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.



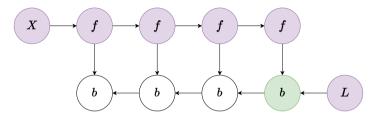


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• All activations f are kept in memory after the forward pass.

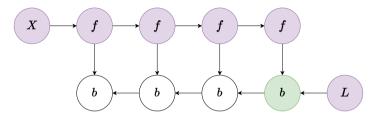


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• All activations f are kept in memory after the forward pass.

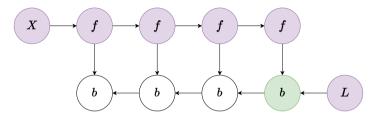


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- All activations f are kept in memory after the forward pass.
 - Optimal in terms of computation: it only computes each node once.

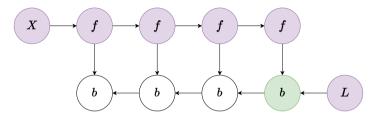


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- All activations f are kept in memory after the forward pass.
 - Optimal in terms of computation: it only computes each node once.

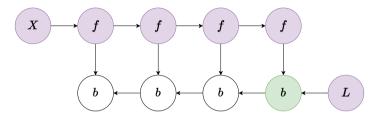
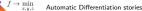


Figure 12: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- All activations f are kept in memory after the forward pass.
 - Optimal in terms of computation: it only computes each node once.

• High memory usage. The memory usage grows linearly with the number of layers in the neural network.



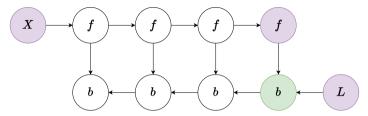


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

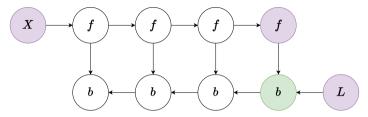


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• Each activation *f* is recalculated as needed.

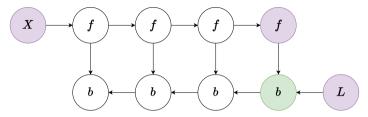


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• Each activation *f* is recalculated as needed.

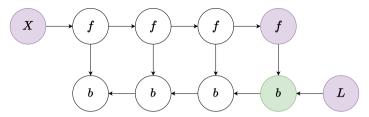


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- Each activation *f* is recalculated as needed.
 - Optimal in terms of memory: there is no need to store all activations in memory.

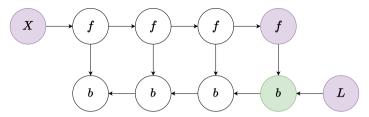


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- Each activation *f* is recalculated as needed.
 - Optimal in terms of memory: there is no need to store all activations in memory.

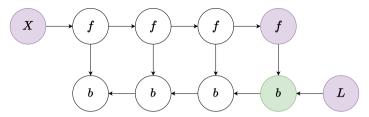


Figure 13: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- Each activation *f* is recalculated as needed.
 - Optimal in terms of memory: there is no need to store all activations in memory.

• Computationally inefficient. The number of node evaluations scales with n^2 , whereas it vanilla backprop scaled as n: each of the n nodes is recomputed on the order of n times.

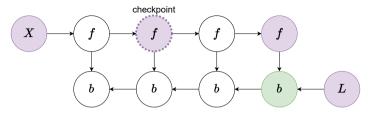


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.



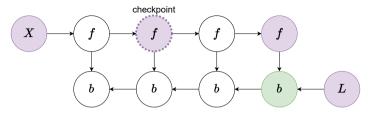


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• Trade-off between the **vanilla** and **memory poor** approaches. The strategy is to mark a subset of the neural net activations as checkpoint nodes, that will be stored in memory.

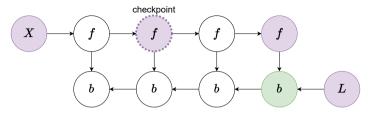


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

• Trade-off between the **vanilla** and **memory poor** approaches. The strategy is to mark a subset of the neural net activations as checkpoint nodes, that will be stored in memory.

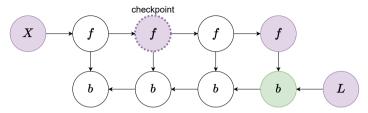


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

 Trade-off between the vanilla and memory poor approaches. The strategy is to mark a subset of the neural net activations as checkpoint nodes, that will be stored in memory.

• Faster recalculation of activations f. We only need to recompute the nodes between a b node and the last checkpoint preceding it when computing that b node during backprop.



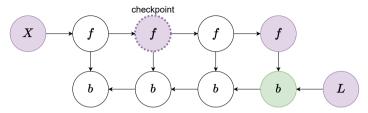


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

 Trade-off between the vanilla and memory poor approaches. The strategy is to mark a subset of the neural net activations as checkpoint nodes, that will be stored in memory.

• Faster recalculation of activations f. We only need to recompute the nodes between a b node and the last checkpoint preceding it when computing that b node during backprop.



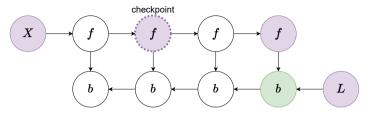


Figure 14: Computation graph for obtaining gradients for a simple feed-forward neural network with n layers. The purple color indicates nodes that are stored in memory.

- Trade-off between the vanilla and memory poor approaches. The strategy is to mark a subset of the neural net activations as checkpoint nodes, that will be stored in memory.
 - Faster recalculation of activations f. We only need to recompute the nodes between a b node and the last checkpoint preceding it when computing that b node during backprop.
 - Memory consumption depends on the number of checkpoints. More effective then vanilla approach.

Gradient checkpointing visualization

The animated visualization of the above approaches \mathbf{O} An example of using a gradient checkpointing \mathbf{O}

Large batch training

