

# Fundamentals of Machine Learning:

## Introduction to Deep Learning

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Prof. Andrew D. Bagdanov (`andrew.bagdanov AT unifi.it`)



UNIVERSITÀ  
DEGLI STUDI  
FIRENZE

**DINFO**  
DIPARTIMENTO DI  
INGEGNERIA  
DELL'INFORMAZIONE

Introduction

Connectionism: The Old School

Connectionism: The New School

Building and Training Deep Networks

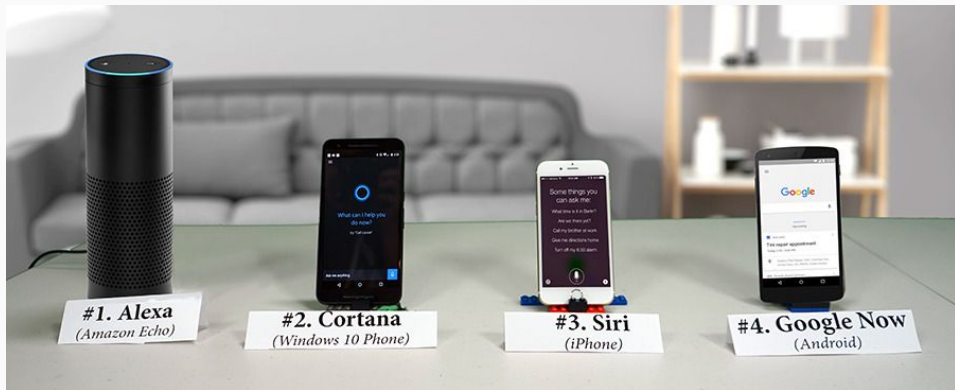
Discussion

# Introduction

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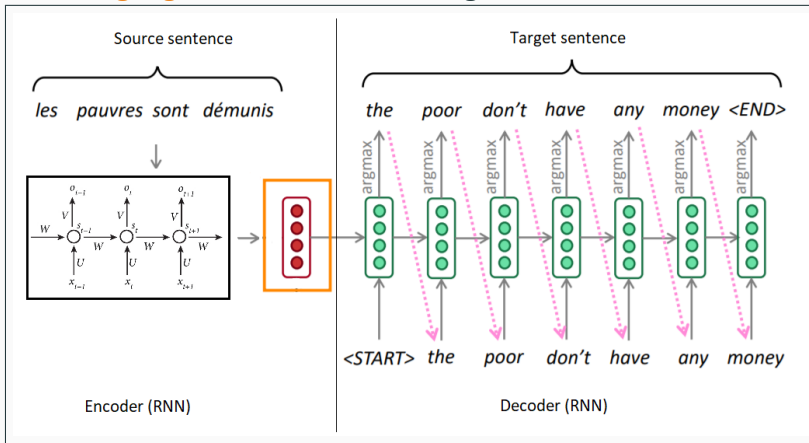
# Digital assistants

- Deep learning is **profoundly** changing our lives.



# Natural language processing

- Deep Recurrent Neural Networks (RNNs) are powering the latest generation of natural language translation technologies.





# Image captioning

- Convolutional Neural Networks (CNNs) are able to extract high-level semantics from images.


Train


**COCO Captions: 80 Classes**


 Two pug **dogs** sitting on a **bench** at the beach.


 A **child** is sitting on a **couch** and holding an **umbrella**.


**Open Images: 600 Classes**


 **Goat**

 **Artichoke**

 **Accordion**


 **Dolphin**

 **Waffle**


 **Balloon**

nocaps Val / Test


**In-Domain: Only COCO Classes**

 The **person** in the brown suit is directing a **dog**.

**Near-Domain: COCO & Novel Classes**

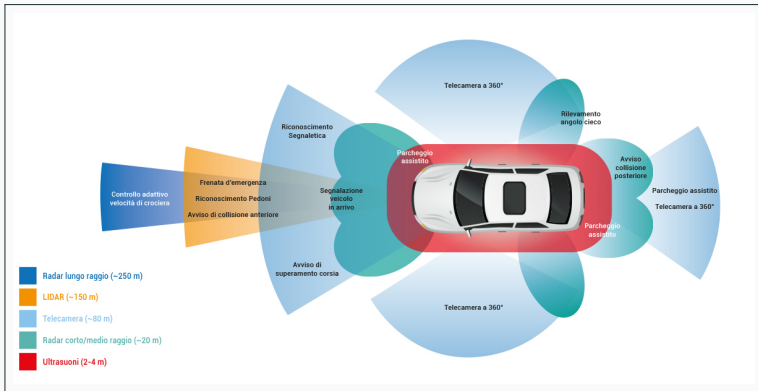
 A **person** holding a black **umbrella** and an **accordion**.

**Out-of-Domain: Only Novel Classes**

 Some **dolphins** are swimming close to the base of the ocean.

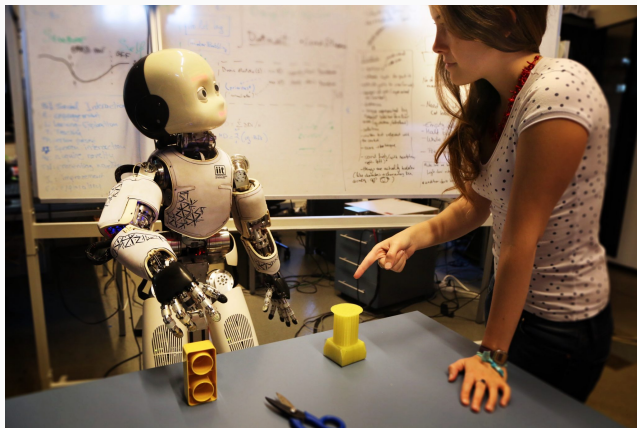
# Self-driving cars

- CNNs are able to integrate multi-modal inputs and are driving the latest advances in Automatic Driving Assistance (ADAS) systems.



# Reinforcement learning

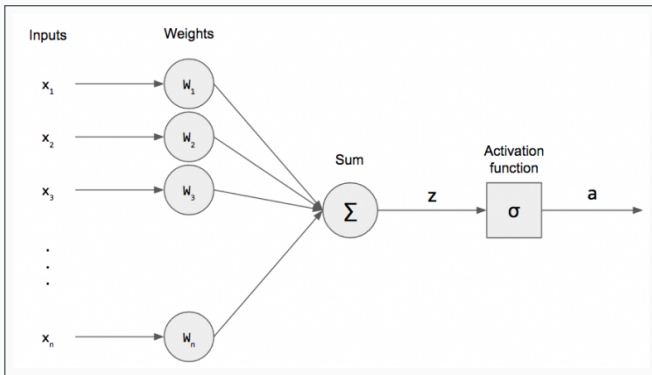
- Deep Reinforcement Learning is being used to train robots who can learn from experience and interactions with their environment.





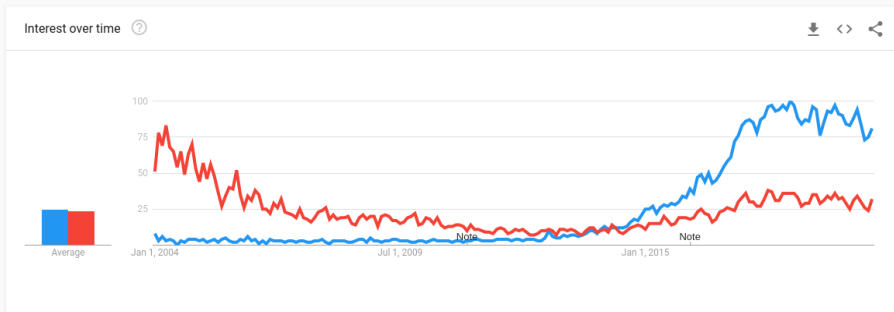
# All thanks to...

- The humble Neural Network.
- Artificial Neural Networks (ANNs) are extremely simple, yet also extremely powerful models.
- They are, in fact, universal function approximators.



# Neural Networks are not new

- As we will see, **neural networks** have a storied history.
- **Deep Learning**, however, is their modern incarnation.



# Overview

- Today we will see what puts the **deep** into **Deep Learning**.
- We will start with an overview of some **historical milestones** in the development of artificial neural networks.
- Then we will look at how modern deep neural networks are **actually built**:
- We will see how the basic **Multilayer Perceptron (MLP)** model provides a **modular** architecture for machine learning problems.
- And we will see how modern tools (e.g. **PyTorch**) makes it easy to apply Deep Models to new problems.

# Lecture objectives

After this lecture you will:

- Understand what **connectionist models** are and how they allow us to compute nonlinear functions of inputs.
- Understand how the **perceptron** works and how its parameters are estimated via **error correction**.
- Understand how the **multi-layer perceptron** generalized the perceptron and how it uses **input**, **hidden**, and **output** layers to represent feedforward computations of inputs to outputs.

## Connectionism: The Old School

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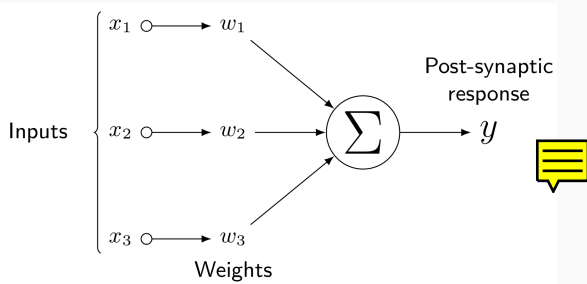
# What is a Deep Neural Network?

- Neural Networks are connectionist models.
- Connectionism has deep roots reaching back to Classical Greece.
- To understand this rich inheritance it is useful to go back in time and trace the roots of modern Deep Models.
- Connectionism arose from the neuroscience and psychological research communities of the 1940s and 1950s.
- These were the nascent beginning of what would become Cognitive Science.
- Though founded on solid experimental practice, what was lacking was any sort of computation basis for learning.

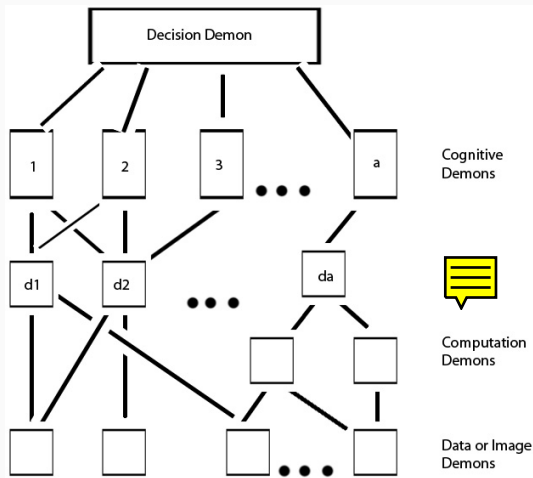
# Connectionism: Hebbian Learning

- One of the first **concrete** learning rules for connectionist models (both artificial and biological).
- **Hebb's Rule**: if cell A consistently contributes to the activity of cell B, then the synapse from A to B should be strengthened.
- **More quaintly**: *neurons that fire together, wire together; neurons that fire out of sync, fail to link.*

$$\begin{aligned}W_{ij} &= x_i x_j \\W_{ij} &= \frac{1}{N} \sum_p x_i^p x_j^p \\ \Delta W_i &= \eta x_i y \\ &= \eta x_i \sum_j W_j x_j\end{aligned}$$



# Connectionism: The Pandemonium Model



- In 1958 Selfridge proposed a multi-layer, **parallel** model of machine learning.
- The model consists of four layers, each inhabited by **demons**.
- Network architecture fixed *a priori*, connections updated using **supervised** learning.
- Demons **yell upwards**, higher-level ones listen and respond.
- **High-worth** demons can replace low-worth ones via **combination**.



# Connectionism: The Perceptron



- The **Perceptron** is probably the simplest (and most famous) feedforward neural network.
- The **perceptron algorithm** was invented by Rosenblatt in 1958.
- It was designed to be a **machine**, and its original purpose was to perform **image recognition**.

$$f(x) = \begin{cases} 1 & \text{if } x > 0 \\ 0 & \text{otherwise} \end{cases}$$



---

The perceptron algorithm

---

**Input:**  $D = \{ (x_i, y_i) \}_{i=1}^N$  (training data)

**Output:** learned weights  $\mathbf{w}$

$\mathbf{w}_0 \leftarrow$  random initialization

$t \leftarrow 1$

**while** not converged **do**

**for**  $(\mathbf{x}, y) \in D$  **do**

$$\hat{y} = f(\mathbf{w}^T \mathbf{x})$$

$$\mathbf{w}_t \leftarrow \mathbf{w}_{t-1} + \eta(y - \hat{y})\mathbf{x}$$

$$t \leftarrow t + 1$$

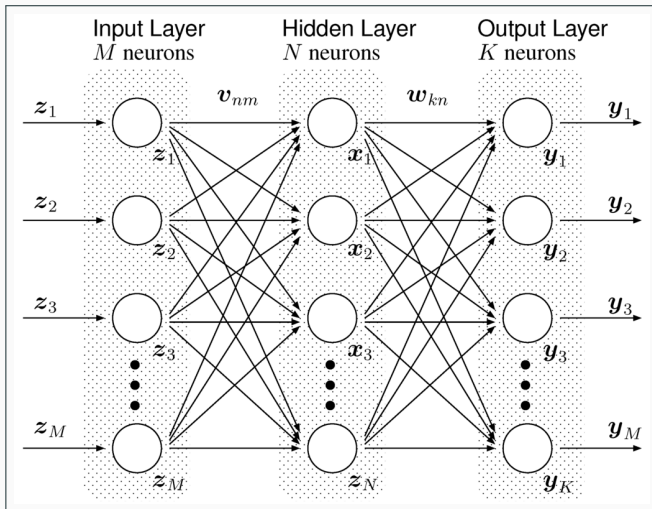


## Connectionism: The New School

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# The Multilayer Perceptron

- Let's look at a simple **Neural Network** architecture known as the **Multilayer Perceptron (MLP)**:



# The Multilayer Perceptron (continued)

- The MLP equation (one hidden layer):

$$\hat{y}(x) = \sigma(w_2^T \sigma(w_1^T x + b_1) + b_2)$$



- Except for the activation function  $\sigma$ , this is a linear system.
- Common activation functions (elementwise):
  - $\sigma(x) = \tanh(x)$
  - $\sigma(x) = (1 + e^{-x})^{-1}$
  - $\sigma(x) = \frac{\exp(x)}{\sum_i \exp(x_i)}$  (softmax, used for outputs).



# Training an MLP

- How do you train a model?
- Decide on a **loss function** (like the negative log-likelihood):

$$L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x})) = -\frac{1}{C} \sum_i y_i \log(\hat{y}_i)$$

- And perform **gradient descent** w.r.t. **all model parameters**:

$$\boldsymbol{\theta}_{n+1} = \boldsymbol{\theta}_n - \varepsilon \nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}))$$

$$\boldsymbol{\theta}_{n+1} = \boldsymbol{\theta}_n - \varepsilon \frac{1}{N} \sum_{i=1}^N \nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}_i))$$



- Where  $\varepsilon$  is the **learning rate**.
- The standard algorithm for this is known as **backpropagation** and it is very clever and efficient.

## Training an MLP (continued)

- OK, we perform **gradient descent** like this:

$$\boldsymbol{\theta}_{n+1} = \boldsymbol{\theta}_n - \varepsilon \frac{1}{N} \sum_{i=1}^N \nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}_i))$$

## Training an MLP (continued)

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- But... How do I compute that gradient  $\nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}_i))$  for **non-trivial** models?

## Training an MLP (continued)

- OK, we perform **gradient descent** like this:

$$\boldsymbol{\theta}_{n+1} = \boldsymbol{\theta}_n - \varepsilon \frac{1}{N} \sum_{i=1}^N \nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}_i))$$

- But... How do I compute that gradient  $\nabla_{\boldsymbol{\theta}} L(\mathbf{y}, \hat{\mathbf{y}}(\mathbf{x}_i))$  for **non-trivial** models?
- The standard algorithm for this is known as **backpropagation** and it is very clever and efficient.
- But... We will defer a **detailed** discussion of backpropagation for the next lecture.



# Stochastic Gradient Descent (SGD)

- **Problem:** what happens if  $N$  (the number of training samples) is **very large**?
- Well, we end up taking **very** slow steps – each iteration of gradient descent is an **average** over the entire dataset.
- **Solution:** approximate the **true** gradient with the gradient at a **single** training example:


## Online Stochastic Gradient Descent



- Choose an initial vector of parameters  $\theta$  and learning rate  $\eta$ .
- Repeat until an approximate **minimum** is found:
  1. **Randomly shuffle training samples** in  $D$ .
  2. For  $(\mathbf{x}, y) \in D$ :

$$\theta := \theta - \eta \nabla_{\theta} \mathcal{L}(\{\mathbf{x}, y\}; \theta)$$

## Stochastic Gradient Descent (continued)

- **Another problem**: evaluating the gradient on **single** examples leads to **very noisy** steps in parameter space.
- One trick to mitigate this is to use **momentum**: keep a running average of gradients that is **slowly** updated.
- Another solution is to use **mini-batches**: instead of a single sample, average the gradients over a **small batch** of samples. 
- It is common to use a combination of **mini-batches** and **momentum** to stabilize training.

# SGD Terminology

- Some useful terminology for deep learning optimization:
  - 1 epoch: one complete pass over the data.
  - 1 iteration: a single gradient step.
  - $N$ : number of training samples.
  - $B$ : batch size.



Algorithm	iterations per epoch
Batch gradient descent	1
Stochastic Gradient Descent	$N$
Mini-batch Gradient Descent	$\frac{N}{B}$

# Building and Training Deep Networks

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# The ingredients

- **Network Definition**: A neural network **model** must be defined that precisely describes the transformation from input to output (e.g. a Multi-layer Perceptron with one hidden layer). The model is typically – but not necessarily – defined in terms of pre-defined, composable **modules**.
- **Dataset and DataLoader**: To train a model, we need some **data** (duh). Data management is **critical** in Deep Learning, and a **DataLoader** is responsible for loading, transforming, and **batching** data for training and inference.
- **A Training Loop**: Neural Networks are not **black boxes**, and their training can be **delicate** and **subtle**. A **training loop** applies an iterative optimization algorithm over **training batches**.



## Defining a network: Basics

- A **network architecture** defines the family of functions we use as **models**.
- It explicitly defines the **parameterized** family of functions  $\mathcal{H}$  we are searching over.
- An architecture is a **Directed Acyclic Graph (DAG)** defining the connections between basic **computational blocks**.
- We call these blocks **layers** and they are reusable **building blocks**.

# Defining a network: The PyTorch Tensor

- **Tensors** are a specialized data structure similar to arrays and matrices.
- In **PyTorch**, tensors are used to:
  - Encode the **inputs** and **outputs** of a model.
  - Encode the **parameters** of a model.
- Tensors vs. NumPy **ndarrays**:
  - Tensors can run on **GPUs or other hardware accelerators.**
  - Tensors and NumPy arrays can share the same underlying memory, eliminating data copying.
- **Tensors are optimized for automatic differentiation.**



# Defining a network: Tensor Operations

- PyTorch tensors support over 100 tensor operations for:
  - Arithmetic
  - Linear algebra
  - Matrix manipulation (transposing, indexing, slicing)
  - Sampling
  - And much, much more.
- Operations can be run on GPU (typically faster than CPU)
  - Default tensor creation on CPU
  - Move to GPU using .to method (check GPU availability first)



## Defining a network: The **tedious** way

- We **can** define our network model completely using low-level tensors.
- For **linear regression** it might look **something** like this:

```
# Define our model parameters, initialize randomly.  
w = torch.randn(Xs.shape[1])  
b = torch.randn(1)  
  
# Try it out by predicting on training set.  
Xs_tr @ w + b  
  
--> tensor([-428.8886, -384.7284, -212.2810, ..., -286.6152, -326.9498, -162.2029])
```

## Defining a network: The tedious way

- OK, but that just looks like NumPy with extra steps.
- The real magic is in the construction of the computational graph.
- And what we can do with the computational graph.

```
# A quadratic loss function.
```

```
def l2loss(y, y_hat):  
    return ((y - y_hat) ** 2.0).mean()
```

```
# Define our model parameters, initialize randomly.
```

```
w = torch.randn(Xs.shape[1], requires_grad=True)
```

```
b = torch.randn(1, requires_grad=True)
```

```
# Compute the loss on the training set.
```

```
l2loss(ys_tr, Xs_tr @ w + b)
```

```
--> tensor(909343.6875, grad_fn=<MeanBackward0>)
```

## Defining a network: The **tedious** way

- Our computational graph is **instrumented** to compute gradients!
- Gradients of the loss wrt tensors with **requires\_grad=True** will be computed:

```
# Compute the gradient of the loss function.
```

```
loss = l2loss(ys_tr, Xs_tr @ w + b)
```

```
loss.backward()
```

```
print(f'Gradient wrt w: {w.grad}')
```

```
print(f'Gradient wrt b: {b.grad}')
```

```
-->
```

```
Gradient wrt w: tensor([-6.7520e+03, -4.5452e+04, -9.1817e+03, -1.8607e+03,  
                        -4.0045e+06, -6.9095e+03, -6.1650e+04,  2.0769e+05])
```

```
Gradient wrt b: tensor([-1739.1476])
```

## Defining a network: The **tedious** way

- We **could** perform gradient descent, but everything must be done **just right**:

```
# Define our model parameters, initialize randomly.  
w = torch.normal(0.0, np.sqrt(1.0/((8+1)/2)), size=(Xs.shape[1],)).requires_grad_()  
b = torch.tensor(1.0, requires_grad=True)  
  
# Standardize dataset.  
Xs_tr -= Xs_tr.mean(0)  
Xs_tr /= Xs_tr.std(0)
```

## Defining a network: The **tedious** way

- We **could** perform gradient descent, but everything must be done **just right**:

```
# Training hyperparameters.
```

```
epochs = 1000
```

```
lr = 1e-2
```

```
# Training loop.
```

```
losses = []
```

```
for it in range(epochs):
```

```
    loss = l2loss(ys_tr, Xs_tr @ w + b)
```

```
    loss.backward()
```

```
    w.data -= lr * w.grad.data
```

```
    b.data -= lr * b.grad.data
```

```
    w.grad.data.zero_()
```

```
    b.grad.data.zero_()
```

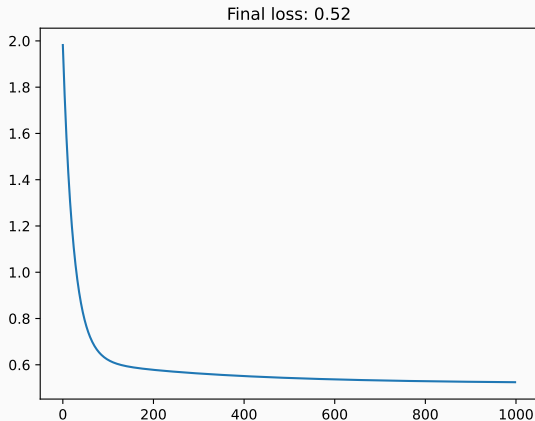
```
    losses.append(loss.item())
```

```
# Plot losses and report last loss (MSE).
```

```
plt.plot(losses[1:]); plt.title(f'Final loss: {losses[-1]}')
```

## Defining a network: The **tedious** way

- We **could** perform gradient descent, but everything must be done **just right**:



## Defining a network: The **modular** way

- OK, that **really** sucks. And is **error-prone**. And doesn't **scale**. And has lots of **other problems**.
- Instead, we will make use of the **reusable modules** from the **torch.nn** package.
- This **PyTorch namespace** contains implementations of **very many** types of layers. **Let's take a look...**
- The simplest way to define out **linear regression model** is:

```
model = nn.Linear(8, 1)
print(list(model.parameters()))
--> [Parameter containing:
      tensor([[-0.1241,  0.2113, -0.1457, -0.2680, -0.3201, -0.2886, -0.2482, -0.0639]],
            requires_grad=True),
      Parameter containing: tensor([0.3110], requires_grad=True)]
```

## Defining a network: The **modular** way

- The classes from the `torch.nn` package **encapsulate** all parameters and provide a uniform **API** to access them.
- Their **instances** act like **functions**:

```
# Our linear model.  
model = nn.Linear(8, 1)  
  
# Training loop.  
losses = []  
for it in range(epochs):  
    loss = l2loss(ys_tr, model(Xs_tr).flatten())  
    model.zero_grad()  
    loss.backward()  
    for p in model.parameters():  
        p.data -= lr * p.grad.data  
    losses.append(loss.item())
```



## Defining a network: The **modular** way

- The **modularity** really shines when layers are **composed**:

```
# Our MLP model.
```

```
model = nn.Sequential(nn.Linear(8, 16), nn.Tanh(), nn.Linear(16, 1))
```

```
# Training loop.
```

```
losses_mod = []
```

```
for it in range(epochs):
```

```
    loss = l2loss(ys_tr, model(Xs_tr).flatten())
```

```
    model.zero_grad()
```

```
    loss.backward()
```

```
    for p in model.parameters():
```

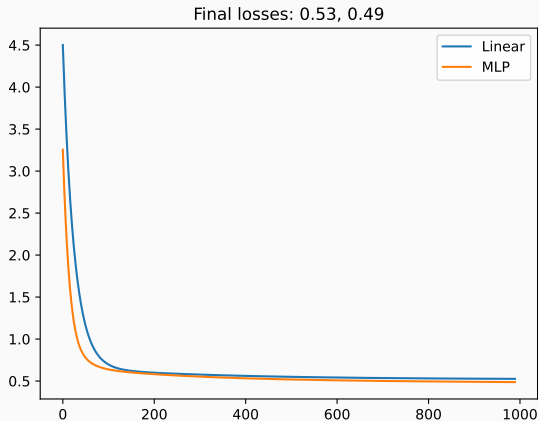
```
        p.data -= lr * p.grad.data
```

```
    losses_mod.append(loss.item())
```



## Defining a network: The **modular** way

- Which makes **model selection** reasonably agile:



## Defining a network: The **modular** way

- The most **general** way to define models is to **extend** `nn.Module`.
- This way we have **complete access** to everything and **control** over composition:

```
class LinRegMLP(nn.Module):
    def __init__(self, num_in, num_hidden, activation=F.tanh):
        super().__init__()
        self.activation = activation
        self.fc1 = nn.Linear(num_in, num_hidden)
        self.out = nn.Linear(num_hidden, 1)

    def forward(self, xs):
        return self.out(self.activation(self.fc1(xs)))
```

## Dataset and DataLoader: Basics

- Data is at the **heart** of Deep Learning.
- The important breakthroughs in Deep Learning would not have been possible without **massive** datasets.
- **Annotated** or **unlabeled** data may power training of Deep Models.
- Model architectures are **strongly** related to the **type** of data.
- Dataset **size** and **annotation quality** will determine how large of a model we can train.

# Dataset and DataLoader: Batching

- In the examples above we are performing **batch gradient descent**.
- Usually this is **inadvisable**, so we use a **DataLoader** which is a **batched iterator** over a dataset:

```
from torch.utils.data import DataLoader, TensorDataset

# Create a TensorDataset
ds_train = TensorDataset(Xs_tr, ys_tr)

# Create a DataLoader
dl_train = DataLoader(ds_train, batch_size=512, shuffle=True)

for (Xs, ys) in dl_train:
    print(f'Xs: {Xs.shape}, ys: {ys.shape}')
-->
Xs: torch.Size([512, 8]), ys: torch.Size([512])
Xs: torch.Size([512, 8]), ys: torch.Size([512])
...
Xs: torch.Size([120, 8]), ys: torch.Size([120])
```



## Dataset and DataLoader: The rest of the story

- Datasets for Deep Learning tend to be **massive**, and data loading is a **critical** issue.
- The **DataLoader** might be responsible for **loading** data from disk, **transforming it** (e.g. **standardization** or **data augmentation**), and **batching** it for training and validation.

# Optimization: Training Deep Models

- Without **optimization** we cannot get anything out of data.
- There are a variety of **optimization algorithms** that manage training of Deep networks.
- 99% of them are based on **gradient-based optimization** (i.e. Stochastic Gradient Descent).
- There is also a (rather large) **bag of tricks** needed to avoid common training pitfalls.

## Optimization: Training Deep Models

- The `torch.nn` module provides classes to encapsulate details of model definitions.
- In the `torch.optim` package you will find implementations of several optimization algorithms that encapsulate the details of model training:
  - `torch.optim.SGD`: Implements the standard Stochastic Gradient Descent (SGD) algorithm, with optional momentum.
  - `torch.optim.Adam`: Implements the Adaptive Moment Estimation (Adam) algorithm, which augments SGD with adaptive, per-parameter learning rates.



# Optimization: Putting Everything Together

- First some **imports** and a **model definition**:

```
import torch.nn as nn
import torch.nn.functional as F
from torch.optim import SGD, Adam
from tqdm.notebook import tqdm

# Define our model.
class LinRegMLP(nn.Module):
    def __init__(self, num_in, num_hidden, activation=F.tanh):
        super().__init__()
        self.activation = activation
        self.fc1 = nn.Linear(num_in, num_hidden)
        self.out = nn.Linear(num_hidden, 1)

    def forward(self, xs):
        return self.out(self.activation(self.fc1(xs)))
```

# Optimization: Putting Everything Together

- And a function to **train** the model:

```
def train(model, dl_train, dl_val=None, epochs=100, lr=1e-3):
    opt = Adam(model.parameters(), lr=lr) # We'll use Adam.
    (train_losses, val_losses) = ([], [])

    # Iterate for required number of epochs.
    for epoch in tqdm(range(epochs)):
        # Iterate over all batches.
        epoch_loss = 0.0
        for (Xs, ys) in dl_train:
            ys_hat = model(Xs).flatten() # Forward pass and loss computation.
            loss = F.mse_loss(ys, ys_hat)
            opt.zero_grad() # Take a gradient step.
            loss.backward()
            opt.step()
            epoch_loss += loss.item()
        train_losses.append(epoch_loss / len(dl_train)) # Collect average epoch loss.

    # If we have a validation dataloader, compute average loss over it.
    if dl_val:
        with torch.no_grad():
            val_losses.append(np.mean([F.mse_loss(ys, model(Xs).flatten()).item() for (Xs, ys) in dl_val]))
    return (train_losses, val_losses)
```

# Optimization: Putting Everything Together

- And (finally) we can **use** this training loop:

```
# Training hyperparameters.
```

```
lr = 1e-3
```

```
epochs = 1500
```

```
batch_size = 512
```

```
inner_size = 16
```

```
# Create datasets and DataLoaders.
```

```
ds_train = TensorDataset(Xs_tr, ys_tr)
```



```
ds_val = TensorDataset(Xs_te, ys_te)
```

```
dl_train = DataLoader(ds_train, batch_size=batch_size, shuffle=True)
```

```
dl_val = DataLoader(ds_val, batch_size=batch_size, shuffle=False)
```



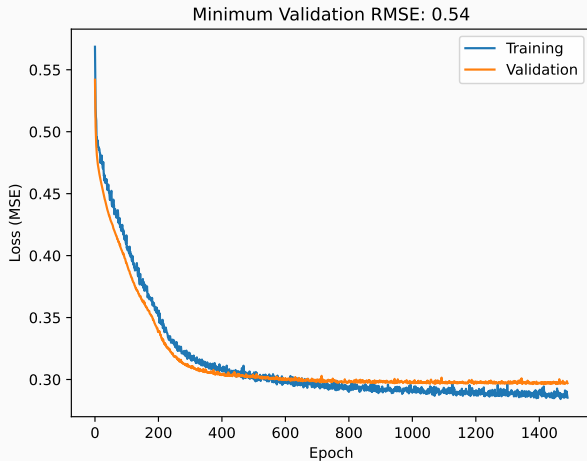
```
# Instantiate model, train it, and plot losses.
```

```
model = LinRegMLP(8, inner_size)
```

```
(train_losses, val_losses) = train(model, dl_train, dl_val, lr=lr, epochs=epochs)
```

# Optimization: Putting Everything Together

- And arrive at something like this:



## Discussion

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## Connectionist models and the state-of-the-art

- **Deep Neural Networks** are connectionist models that represent the **state-of-the-art** for many learning problems today.
- They are usually trained with **backpropagation**.
- Since they typically have **very many parameters** they must be trained with **extreme care**.
- They also typically require **more training data** than classical approaches in order to generalize.
- Note that there is nothing **fundamentally** different about these models with respect to the **classical** models we have already seen.
- They simply **represent a much richer hypothesis set of functions  $\mathcal{H}$  over which we search.**

# The Embarrassment of Choice

- **PyTorch** is the **de facto** framework of choice for most Deep Learning projects.
- It is not the **only one**:
  - **Tensorflow/Keras**: a graph-based, automatic differentiating framework for deep learning in Python – now with **JAX** and **PyTorch** backends!
  - **JAX**: A Python library for accelerator-oriented array computation and program transformation – sold as **differentiable Python programming**.
- Moreover, there are many **frameworks** for tracking experiments:
  - **Tensorboard**: A suite of web applications for inspecting and understanding your TensorFlow (and PyTorch, and JAX, and Keras, and...) runs and graphs.
  - **Weights and Biases**: Track, version and visualize with just 5 lines of code, reproduce any model checkpoints, monitor CPU and GPU usage in real time.
  - **Comet**: Provides an end-to-end model evaluation platform for AI developers, with best in class LLM evaluations, experiment tracking, and production monitoring.

# Reading and Homework Assignments

## Reading Assignment:

- Bishop: Chapter 5
- PyTorch: [This is an excellent introductory tutorial](#)