L12 – Statistical Mechanics 4

Machine learning

Neural networks

What is a Neural Network?

Adapted from: <u>www.3blue1brown.com/lessons/neural-networks</u>

Plain vanilla (aka "multilayer perceptron")









 0
 0.7
 0.00
 0.0
 0

 0
 1.0
 0.6
 0.1
 0

 1
 1.0
 1.0
 0.2
 0

0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.2 0.5 0.9 0.9 1.0 1.0 1.0 1.0 1.0 1.0 1.0 0.6 0.1 0.0 0.0 0.0 0.0 0.0 0.0 1.0 1.0 0.9 0.5 0.5 0.5 0.5 0.7 1.0 1.0 1.0 1.0 2.2 0.0 0.1 0.8 0.8 0.8 1.0 1.0 1.0 1.0 0.9 0.4 0.1 0.0 0.0 0.0 0.0 0.0 0.0 0.0 0.0



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28

$28 \times 28 = 784$



... 00000000 F $\not\models$ $\begin{array}{c} 0 \\ 0 \\ 1 \\ 2 \\ 3 \\ 0 \\ 4 \\ 5 \\ 0 \\ 6 \\ 7 \\ 8 \\ 9 \\ \end{array}$ 0000000 \varkappa X Ľ E

784



0000000 ··· 00000000 F $\not\models$ $egin{array}{c} 0 \\ 0 \\ 1 \\ 0 \\ 2 \\ 3 \\ 0 \\ 4 \\ 0 \\ 5 \\ 0 \\ 6 \\ 0 \\ 7 \\ 0 \\ 8 \\ 0 \\ 9 \end{array}$ \geq

784



$\sigma(w_1a_1 + w_2a_2 + w_3a_3 + \dots + w_na_n-10)$ "bias"

Only activate meaningfully when weighted sum > 10



 $784 \times 16 + 16 \times 16 + 16 \times 10$ weights

 $\begin{array}{c} 16 + 16 + 10 \\ \text{biases} \end{array}$

 $13,\!002$

784

Subscript corresponds to a neuron in the layer

Sigmoid

$$= \sigma \left(w_{0,0} \ a_{0}^{(0)} + w_{0,1} \ a_{1}^{(0)} + \dots + w_{0,n} \ a_{n}^{(0)} + b_{0} \right)$$
Bias

$$\frac{w_{0,0} \ w_{0,1} \ \dots \ w_{0,n}}{w_{1,0} \ w_{1,1} \ \dots \ w_{1,n}} \left[\begin{array}{c} a_{0}^{(0)} \\ a_{1}^{(0)} \\ \end{array} \right]$$

$$\vdots$$
 \vdots \ddots \vdots $u_{k,0}$ $w_{k,1}$ \ldots $w_{k,n}$ $a_n^{(0)}$

Sigmoid

$$\int_{0}^{(1)} = \sigma \left(w_{0,0} \ a_{0}^{(0)} + w_{0,1} \ a_{1}^{(0)} + \dots + w_{0,n} \ a_{n}^{(0)} + b_{0} \right) \\
\overset{(1)}{\underset{(1)}{\overset{(1)}{\underset{($$

Sigmoid

$$a_{0}^{(1)} = \overset{\downarrow}{\sigma} \left(w_{0,0} \ a_{0}^{(0)} + w_{0,1} \ a_{1}^{(0)} + \dots + w_{0,n} \ a_{n}^{(0)} + b_{0} \right)$$
Bias

 $a_0^{(0)}$ $w_{0,0} \ w_{0,1} \ \dots \ w_{0,n}$ b_0 $a_1^{(0)}$ b_1 $w_{1,0} \ w_{1,1} \ \dots \ w_{1,n}$ $a_n^{(0)}$ b_n $w_{k,0}$ $w_{k,1}$ \ldots $w_{k,n}$

Neural network function

10 outputs

A NN with L hidden layers takes an input vector \mathbf{x} and returns an output vector \mathbf{y} though the forward equations chain

$$\mathbf{a}^{(1)} = \sigma \left(\mathbf{w}^{(0)} \mathbf{x} + \mathbf{b}^{(0)} \right)$$
$$\mathbf{a}^{(2)} = \sigma \left(\mathbf{w}^{(1)} \mathbf{a}^{(1)} + \mathbf{b}^{(1)} \right)$$
$$\vdots$$
$$\mathbf{a}^{(L)} = \sigma \left(\mathbf{w}^{(L-1)} \mathbf{a}^{(L-1)} + \mathbf{b}^{(L-1)} \right)$$
$$\mathbf{y} = \sigma \left(\mathbf{w}^{(L)} \mathbf{a}^{(L)} + \mathbf{b}^{(L)} \right)$$

The NN is a function of ${\bf x}$ with a parametric dependence on ${\bf w}$ and ${\bf b}$

$$\mathbf{y} = f(\mathbf{x}; \mathbf{w}, \mathbf{b})$$

The cost of learning

Adapted from: <u>www.3blue1brown.com/lessons/gradient-descent</u>

(0,0)(4,6)(3,3)(4,6)(7,7)(8,8)(0,0)(7,9) $(\mathbf{5}, 5)(\mathbf{4}, 4)(\mathbf{3}, 3)(\mathbf{4}, 6)(\mathbf{5}, 5)(\mathbf{8}, 8)(\mathbf{9}, 9)(\mathbf{5}, 5)$ (4,4)(4,4)(7,7)(2,2)(0,0)(3,3)(2,2)(8,8)([9,9)([1,1)([9,9)([2,2)([2,2)([2,7)([9,9)([4,4) $(\mathbf{8}, 8)(\mathbf{7}, 7)(\mathbf{4}, 4)(\mathbf{7}, 1)(\mathbf{3}, 3)(\mathbf{7}, 1)(\mathbf{5}, 5)(\mathbf{3}, 3)$ $(\mathbf{2}, 2)(\mathbf{3}, 3)(\mathbf{9}, 9)(\mathbf{0}, 0)(\mathbf{9}, 9)(\mathbf{9}, 9)(\mathbf{1}, 1)(\mathbf{5}, 5)$ $(\mathbf{3}, 8)(\mathbf{4}, 4)(\mathbf{1}, 7)(\mathbf{7}, 7)(\mathbf{4}, 4)(\mathbf{4}, 4)(\mathbf{4}, 4)(\mathbf{3}, 2)$ $(\mathbf{0},0)(\mathbf{7},7)(\mathbf{2},2)(\mathbf{4},4)(\mathbf{8},8)(\mathbf{2},2)(\mathbf{6},6)(\mathbf{9},9)$ ([9,9)([2,2)([5,8)]([2,7)([6,6)]([1,1)([1,1)([2,2))))) $(\mathbf{3},3)(\mathbf{9},9)(\mathbf{1},1)(\mathbf{6},6)(\mathbf{5},5)(\mathbf{1},1)(\mathbf{1},1)(\mathbf{0},0)$

Test on these

What's the "loss" of this difference?

Utter trash

Loss of

$$\begin{array}{l} (0.43-0.00)^2+\\ (0.28-0.00)^2+\\ (0.19-0.00)^2+\\ (0.88-1.00)^2+\\ (0.72-0.00)^2+\\ (0.01-0.00)^2+\\ (0.64-0.00)^2+\\ (0.86-0.00)^2+\\ (0.99-0.00)^2+\\ (0.63-0.00)^2\end{array}$$

Cost function

 $(\textcircled{3},9)(\textcircled{0},0)(\textcircled{2},2)(\textcircled{6},6) \\ (\textcircled{0},0)(\cancel{4},4)(\cancel{6},6)(\cancel{7},7) \\ (\fbox{7},7)(\cancel{8},8)(\cancel{3},3)(\cancel{1},1) \\ (\fbox{1},1)(\cancel{1},1)(\cancel{6},6)(\cancel{3},3) \\ (\fbox{1},1)(\cancel{1},1)(\textcircled{0},0)(\cancel{4},4) \end{cases}$

Lots of training data

3.37

One number

13,002 weights and biases For each training data k, with true value $\hat{\mathbf{y}}_k$ get the **loss function** over the output neurons

$$\mathcal{L}_{k}\left(\mathbf{w},\mathbf{b}\right) = \frac{1}{N_{out}} \sum_{i=1}^{N_{out}} \left(NN\left(x_{k,i};\mathbf{w},\mathbf{b}\right) - \hat{y}_{k,i}\right)^{2}$$

Considering all training data, get the **cost function**

$$C(\mathbf{w},\mathbf{b}) = \frac{1}{N_{train}} \sum_{k=1}^{N_{train}} \mathcal{L}_k(\mathbf{w},\mathbf{b})$$

The optimized NN is the one with **w** and **b** that minimizes *C*(**w**, **b**)

$$\min_{\mathbf{w},\mathbf{b}} C(\mathbf{w},\mathbf{b}) \Longrightarrow \nabla_{\mathbf{w},\mathbf{b}} C(\mathbf{w},\mathbf{b}) = 0$$

How to Train Your Neural Network

Adapted from: <u>www.3blue1brown.com/lessons/gradient-descent</u>

$$-\nabla C(\vec{\mathbf{W}}) = \begin{bmatrix} 0.31 \\ 0.03 \\ -1.25 \\ \vdots \\ 0.78 \\ -0.37 \\ 0.16 \end{bmatrix}$$

3

8

 w_0 should increase somewhat w_1 should increase a little w_2 should decrease a lot

 $\overline{w}_{13,000}$ should increase a lot $w_{13,001}$ should decrease somewhat $w_{13,002}$ should increase a little

The minimization is done with **gradient descent**:

$$\mathbf{w}_{m+1} = \mathbf{w}_m - \gamma_m \frac{\partial C(\mathbf{w}_m, \mathbf{b}_m)}{\partial \mathbf{w}_m} \qquad (\gamma_m - \text{learning rate})$$
$$\mathbf{b}_{m+1} = \mathbf{b}_m - \gamma_m \frac{\partial C(\mathbf{w}_m, \mathbf{b}_m)}{\partial \mathbf{b}_m}$$

Usually, the gradient is computed for a sub-set of **w** and **b** components chosen at random. It is called **stochastic gradient descent**

en.wikipedia.org/wiki/Gradient_descent

Now that we have optimal **w** and **b**, we can use the neural network to identify images that were not in the training set.

This is the basics. But there is so much more...

• Physics-informed neural networks (PINN)

Integrates partial differential equations expressing physical laws into the NN cost function <u>youtu.be/G hlppUWcsc</u>

• Graph neural networks (GNN)

NN for processing data that can be represented as vertices connected by edges <u>distill.pub/2021/gnn-intro/</u>

Convolutional neural networks (CNN)

NN with convolutional layers that capture local patterns and global features in the input data <u>tinyurl.com/convnnet</u>

Generative adversarial networks (GAN)

Two NN – a generator and a discriminator – compete to create realistic-looking outputs <u>tinyurl.com/ganetintro</u>

Transformer neural networks

NN architecture for encoding words, word position, and word's contextual relation with others in the sentence (self-attention). https://youtu.be/zxQyTK8quyY

Artificial inteligence

Machine learning in practice

Machine learning development network

www.v7labs.com/blog/pytorch-vs-tensorflow

Repository Creation Date

www.v7labs.com/blog/pytorch-vs-tensorflow

import torch import torch.nn as nn

```
class SimpleNeuralNetwork(nn.Module):
    def __init__(self):
        super(SimpleNeuralNetwork, self).__init__()
        self.layer1 = nn.Linear(784, 16)
        self.layer2 = nn.Linear(16, 16)
        self.output_layer = nn.Linear(16, 10)
```

```
def forward(self, x):
    x = torch.sigmoid(self.layer1(x))
    x = torch.sigmoid(self.layer2(x))
    x = self.output_layer(x)
    return x
```

Instantiate the model
model = SimpleNeuralNetwork()

<section-header>

Deep Learning With PyTorch

Codemy.com

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In this playlist we'll learn A.I. Deep Learning with Pytorch and Python. If you're new to Artificial Intelligence, this is for you!

tinyurl.com/pytorchlearn

ML for Modeling Materials

AI for theoretical chemistry has been used to

- Search the chemical space of compounds
- Perform dimensionality reduction, clustering, and pattern recognition
- Improve or accelerate quantum chemical methods
- Predict properties as a surrogate approach

Search the chemical space of compounds

GNN-based discovery of new materials

Merchant et al. Nature **2023**. <u>10.1038/s41586-023-06735-9</u>

GNN-based discovery added 381,000 new stable materials to the database

Merchant et al. Nature **2023**. <u>10.1038/s41586-023-06735-9</u>

Perform dimensionality reduction, clustering, and pattern recognition

Hierarchical protocol for the automatic analysis of the ring deformation in surface hopping

Based on

- dimensionality reduction (PCA)
- clustering (DBSCAN + agglomerative clustering)

56.2% C1-puckering Cluster A1B2C2 12.1% C1-puckering Cluster A1B2C3 6.3% C1-puckering NH_2 out-of-plane motion $[+]^a$ Cluster A1B3 7.7% **Cluster A2**

7.5%

Channel

Cytosine

Cluster A1B1 10.2%

Cluster A1B2C1

Zhu et al. PCCP **2022**, 24, 24362

 NH_2 out-of-plane motion $[-]^a$

C1-puckering C10-puckering C=O out-of-plane motion $[-]^a$

Important motion

C1-puckering C=O out-of-plane motion $[+]^a$

Improve or accelerate quantum chemical methods

Density functional from an NN

Input features:

- charge density r,
- norm of charge density
- electron kinetic energy density
- local HF exchange energy densities

"The resulting functional, DM21 (DeepMind 21), correctly describes typical examples of artificial charge delocalization and strong correlation and performs better than traditional functionals on thorough benchmarks for main-group atoms and molecules. DM21 accurately models complex systems such as hydrogen chains, charged DNA base pairs, and diradical transition states."

Kirkpatrick *et al. Science* **2021**, *374*, 1385 Quanta magazine: <u>tinyurl.com/qmdm21</u>

Predict properties as a surrogate approach: ML Potentials

$$\mathbf{R} \rightarrow \left(T_{elec} \left(\mathbf{r} \right) + V \left(\mathbf{r}, \mathbf{R} \right) \right) \varphi \left(\mathbf{r}; \mathbf{R} \right) = \frac{E(\mathbf{R})}{\varphi(\mathbf{r}; \mathbf{R})} \rightarrow E(\mathbf{R})$$

Descriptor

Example: ANI ML Potential

Gao et al. J Chem Inf Model **2020**, 60, 3408

- MD17 Database
- Energy + Force

•
$$N_{train} = 1$$
k; $N_{model} = 20$; $N_{test} = 20$ k

To know more:

3Blue1Brown Course on NN

• <u>www.3blue1brown.com/topics/neural-networks</u>

Kernel Methods

• Pinheiro Jr; Dral, In *Quantum chemistry in the age of machine learning*, **2023**; pp 205

ML Potentials

• Pinheiro Jr *et al*. *Chem Sci* **2021**, *12*, 14396

Deep Learning Applied to Computational Mechanics

• Vu-Quoc; Humer. *Comput Model Eng Sci* **2023**, *137*, 1069

Papers available for download at: <u>amubox.univ-amu.fr/s/xXAiMZrDPb9RMRX</u> (Ask me for the password)

Computational modeling of nanosystems

Epiloge

many

Quantum Statistical Mechanics

Classical Statistical Mechanics

Computational Modeling of Nanosystems

Scientific skills

- Math skills Quantum chemistry
- Molecular dynamics
 - Physical chemistry

• Linear algebra Statistics •

Machine learning

Operational skills

- Modeling
- Programming
- Data processing

large

few

Quantum field theory

Standard model

Quantum **Mechanics**

small

Classical **Mechanics**

> General relativity Dark matter/energy