
tranches Documentation

Release 0.1.0

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Contents

1	User's Guide	3
1.1	Requirements	3
1.2	Issues	3
1.3	Installing tranches	3
2	Machine Learning Basics	5
2.1	The supervised learning problem	5
2.2	Loss and Risk Functions	5
2.3	Gradient Descent	6
2.4	Batch Learning	7
2.5	Metrics	7
3	Neural Networks	9
3.1	Fully connected, feed forward networks	9
3.2	Activation Function	9
3.3	Backpropagation	9
3.4	Initialization	9
3.5	Beyond SGD	9
3.6	Regularization	9
4	The MNIST dataset (yet again)	11
4.1	Model selection	11
4.2	Results	11
5	API Reference	13
5.1	activations	13
5.2	initializers	13
5.3	layers	13
5.4	loss	13
5.5	metrics	13
5.6	network	13
5.7	optimizers	13
5.8	scheduling	13
5.9	validation	13
6	tranches changelog	15
6.1	Version 0.1.1	15

7	License	17
7.1	Authors	17
7.2	Contributing	17
7.3	License Definitions	17
7.4	tranches License	17
	Python Module Index	21

Non-optimised implementation of neural networks for educational purposes.

1.1 Requirements

tranches is build on top of the following libraries:

- *Numpy* (<http://www.numpy.org>)
- *matplotlib* (<http://matplotlib.org/>)

1.2 Issues

Should you encounter any issue with the library you can raise them here: <https://github.com/RonsenbergVI/tranches/issues>

1.3 Installing tranches

1.3.1 Installation with pip (recommended)

The library is available on pypi and to install the last available version, run this command:

```
$ pip3 install tranches
```

To test the installation:

```
>>> import tranches
>>> tranches.__version__
```

this should display the version installed on your system.

1.3.2 Installation from GitHub

tranches releases are also available on github (<https://github.com/ronsensbergVI/tranches>).

You first need to clone (or fork if you want to modify it) and

```
$ git clone https://github.com/ronsensbergVI/tranches.git
$ cd tranches
$ python setup.py build
$ python setup.py install
```

Machine Learning Basics

Deep learning is a vibrant subfield of machine learning.. The field has attracted a lot of attention over the past few years. It is a consequence of the continuous progress made in data collection, processing and storage and of the advances in parallelisation allowing to state of the art models.

Even though deep learning has its specificities, it is based on learning theories and it is worth going over some of the fundamentals. This chapter is a brief introduction to the theory of learning. It is the bare minimum needed to go forward but is in no case complete. Feel free to skip to the next chapter.

2.1 The supervised learning problem

We assume that we have at hand a set of examples $(\mathbf{x}_i, y_i)_{i \in \{0, \dots, N\}}$ where \mathbf{x}_i is a vector of features and y_i are labels that are either draw from a discrete set (classification) or the real line (regression). In other words, we have a set of inputs and their desired output.

The goal of supervised learning is to find a function f that ultimately is able to predict the desired output for new examples by learning from the examples at our disposal. Ideally, to solve this problem, we would like to be able to find the joint probability:

$$F(x, y) = \mathbb{P}(\mathbf{x}, y)$$

The learning problem is usually defined over a parameter set and thus we are looking for a function $f_\theta(x)$ with $x \in \Theta$ and Θ is the parameter set. The learning procedure will define the parameter update rules. The parameters are updated so that the model output $\hat{y}_n = f_\theta(x_n)$ are the closest possible to the expected outputs y_n . We need to define metrics to define how close is the model.

2.2 Loss and Risk Functions

In order to learn the optimal parameters θ^* supervised learning algorithms want to minimise what is referred to as a loss function.

A loss function is defined a map taking as input a pair of elements from the possible of labels to a real number. Ideally we would like to have a function that:

1. is regular enough (continuous and convex is the jackpot but hardly ever the case for any interesting model);
2. sensitive to magnitude (this allows to distinguish between the types of error).

The type of loss function chosen depends on the supervised task. For classification the most common loss function is the cross entropy defined as:

$$C(y, \hat{y}) = \sum_k \mathbb{P}(y = k) \log \mathbb{P}(\hat{y} = k)$$

and for regression it is the mean square root (MSE) defined as:

$$C(y, \hat{y}) = ||y - \hat{y}||$$

The loss function computes the “cost” of the learner’s error for one example (intuitively a zero cost means a correct prediction). We want to minimise the cost for the whole dataset. For this we define the (empirical) risk function as:

$$J(\theta) = \frac{1}{N} \sum_n C(y_n, \hat{y}_n)$$

2.3 Gradient Descent

Gradient based optimisation methods are quite popular in machine learning and particularly in deep learning. The gradient of a function gives meaningful information about how to update its parameters to reduce its value. Gradient based methods are iterative algorithms that produce a sequence (theta_n) of parameters where each new iteration improves the risk function.

1. defining an initial value x_0
2. an iteration process to get x_{n+1} given x_n
3. a convergence criteria to stop the training when reached

The simplest gradient based optimisation method is the gradient (or steepest) descent algorithm where each new iteration is defined by going in small steps in the opposite direction of the risk function gradient:

1. initialise x_0
2. $x_{n+1} = x_n - \gamma \nabla f(x_n)$
3. until $||x_{n+1} - x_n|| < \epsilon$

The interesting feature of this method is that it is guaranteed to reach a local minima for a convex function on a convex set.

The hyperparameter γ is referred to as the learning rate. Choosing a learning rate is no easy task (well nothing is really in deep learning) and is critical for the performance of the training. Choosing a too large learning rate prevents the algorithm to target precisely enough regions of local minima and thus oscillates. A large enough can even prevent the algorithm to converge. Too small a convergence makes the algorithm slow as it takes too many steps towards a local minima.

A common method is to start with a large learning rate when the model starts training and is (arguably) far from the optimal solution and as training goes to decay the rate as the model gets closer to the solution we need to take more precaution and not miss it. The popular learning rate schedules used in practice are:

1. exponential decay:

$$\gamma_r = \gamma_0 \exp(-H \frac{r}{R})$$

2. power decay:

power decay:

$$\gamma_r = \gamma_0(1 - \frac{r}{R})^H)$$

2.4 Batch Learning

2.5 Metrics

CHAPTER 3

Neural Networks

<intro to chapter here>

3.1 Fully connected, feed forward networks

3.2 Activation Function

3.3 Backpropagation

3.4 Initialization

3.5 Beyond SGD

3.6 Regularization

The MNIST dataset (yet again)

<intro to chapter here>

4.1 Model selection

4.2 Results

If you are looking for information on a specific function, class or method, this part of the documentation is for you.

5.1 activations

5.2 initializers

5.3 layers

5.4 loss

5.5 metrics

5.6 network

5.7 optimizers

5.8 scheduling

5.9 validation

CHAPTER 6

tranches changelog

We detail here the changes made to the library

6.1 Version 0.1.1

Release day: September 27 2018

- Initial public release.

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The full license text can be found below (tranches-license).

7.1 Authors

tranches is written and maintained by Rene-Jean Corneille

7.2 Contributing

Contribution will be welcomed once a first stable release is ready.

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Version 2.0, January 2004

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