Recurrent neural networks

A nice subtitle

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Introduction

The model

Def: RNN

Given an input sequence $\{u\}_{t=1,...,T}$, with $u_t \in \mathbb{R}^p$, the output sequence of a RNN $\{y\}_{t=1,...,T}$, with $y_t \in \mathbb{R}^o$, is defined by the following:

$$\mathbf{y}^t \triangleq F(\mathbf{z}^t) \tag{1}$$

$$\mathbf{z}^t \triangleq W^{out} \cdot \mathbf{a}^t + \mathbf{b}^{out} \tag{2}$$

$$\boldsymbol{a}^{t} \triangleq W^{rec} \cdot \boldsymbol{h}^{t-1} + W^{in} \cdot \boldsymbol{u}^{t} + \boldsymbol{b}^{rec}$$
 (3)

$$\boldsymbol{h}^t \triangleq \sigma(\boldsymbol{a}^t) \tag{4}$$

$$\boldsymbol{h}^0 \triangleq \mathbf{0},\tag{5}$$

where $\sigma(\cdot): \mathbb{R} \to \mathbb{R}$ is a non linear function applied element-wise called **activation function**, $F(\cdot)$ is called **output function**.

The parameters of the net are $\{W^{out}, W^{in}, W^{rec}, \boldsymbol{b}^{rec}, \boldsymbol{b}^{out}\}$.

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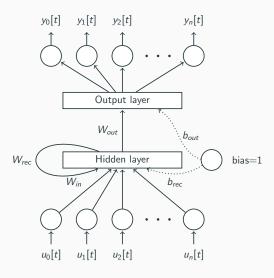


Figure 1: RNN model.

The optimization problem

Given a dataset D:

$$D \triangleq \{\{\overline{\boldsymbol{u}}^{(i)}\}_{t=1,...,T}, \overline{\boldsymbol{u}}_t^{(i)} \in \mathbb{R}^p, \{\overline{\boldsymbol{y}}^{(i)}\}_{t=1,...,T}, \overline{\boldsymbol{y}}_t^{(i)} \in \mathbb{R}^o; i = 1,..., N\}$$
(6)

we define a loss function $L_D(x)$ over D as

$$L_D(\mathbf{x}) \triangleq \frac{1}{N} \sum_{i=1}^{N} \sum_{t=1}^{T} L_t(\overline{\mathbf{y}}_t^{(i)}, \mathbf{y}_t^{(i)}(\mathbf{x})), \tag{7}$$

where $L_t(\cdot, \cdot)$ is an arbitrary loss function for the time step t and x represents all the parameters of the network. The problem is

$$\min_{\mathbf{x}} L_D(\mathbf{x}) \tag{8}$$

Some learning examples

▶ Regression: mean squared error, linear output

$$L(\mathbf{y}, \mathbf{t}) = \frac{1}{M} \sum_{i=1}^{M} (y_i - t_i)^2, \quad F(\mathbf{y}) = \mathbf{y}.$$
 (9)

▶ Binary classification: hinge loss, linear output

$$L(y,t) = \max(0, 1 - t \cdot y), \quad F(y) = y.$$
 (10)

Multi-way classification: cross entropy loss, softmax output

$$L(\mathbf{y}, \mathbf{t}) = -\frac{1}{M} \sum_{i=1}^{M} \log(y_i) \cdot t_i, \quad F(y_j) = \frac{e^{y_j}}{\sum_{i=1}^{M} e^{y_i}}.$$
 (11)

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Stochastic gradient descent (SGD)

Algorithm 1: Stochastic gradient descent

Data:

 $D = \{\langle \boldsymbol{u}^{(i)}, \boldsymbol{y}^{(i)} \rangle\}$: training set

 x_0 : candidate solution

m: size of each mini-batch

Result:

x: solution

- $x \leftarrow x_0$
- 2 while stop criterion do
- 3 $I \leftarrow \text{select } m \text{ training example } \in D$
- 4 $\alpha \leftarrow$ compute learning rate
- 5 $\mathbf{x} \leftarrow \mathbf{x} \alpha \sum_{i \in I} \nabla_{\mathbf{x}} L(\mathbf{x}; \langle \mathbf{u}^{(i)}, \mathbf{y}^{(i)} \rangle)$
- 6 end

Convergence of SDG

- ▶ Nemirovski (2009)[4]: proof of convergence in the convex case
- ▶ there are no theoretical guarantees in the non-convex case
- ▶ in practice it always works: SGD is the standard framework in most of neural networks applications.

The vanishing gradient problem

A pathological problem example

An input sequence:

marker	0	1	0	 0	1	0	0
value	0.3	0.7	0.1	 0.2	0.4	0.6	0.9

The predicted output should be the sum of the two one-marked positions (1.1).

Why is this a difficult problem?

Because of its long time dependencies

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Gradient

Let $L_t(\bar{u}, \bar{y})$ the loss function for the time step t and $g_t(x)$: \mathbb{R} be the function defined by

$$g_t(\mathbf{x}) \triangleq L_t(F(\mathbf{z}^t(\bar{\mathbf{u}};\mathbf{x}),\bar{\mathbf{y}}_t))$$

and

$$g(\mathbf{x}) \triangleq \sum_{t=1}^{T} g_t(\mathbf{x})$$

$$\frac{\partial g}{\partial W^{rec}} = \sum_{t=1}^{T} \nabla L_{t}^{T} \cdot J(F) \cdot \frac{\partial \mathbf{z}^{t}}{\partial \mathbf{a}^{t}} \cdot \frac{\partial \mathbf{a}^{t}}{\partial W^{rec}}$$
(12)

$$= \sum_{t=1}^{T} \frac{\partial g_t}{\partial \mathbf{a}^t} \cdot \frac{\partial \mathbf{a}^t}{\partial W^{rec}}$$
 (13)

Let's consider a single output unit u, and a weight w_{lj} , we have

$$\frac{\partial a_u^t}{\partial w_{lj}} = \sum_{k=1}^t \frac{\partial a_u^t}{\partial a_l^k} \cdot \frac{\partial a_l^k}{\partial w_{lj}} \tag{14}$$

$$=\sum_{k=1}^{t} \delta_{lu}^{tk} \cdot \phi_j^{t-1} \tag{15}$$

where

$$\delta_{lu}^{tk} \triangleq \frac{\partial a_u^t}{\partial a_l^k}.\tag{16}$$

Let P(I) be the set of parents of neuron I, defined as the set of parents in the unfolded network.

$$\delta_{lu}^{tk} = \sum_{h \in P(l)} \delta_{hu}^{tk} \cdot \sigma'(a_h^{t-1}) \cdot w_{hl}$$
 (17)

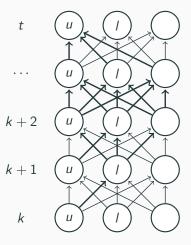


Figure 2: Nodes involved in $\frac{\partial a_u^t}{\partial a_l^t}$.

In matrix notation we have:

$$\frac{\partial \mathbf{a}^{t}}{\partial W^{rec}} = \sum_{k=1}^{t} \frac{\partial \mathbf{a}^{t}}{\partial \mathbf{a}^{k}} \cdot \frac{\partial^{+} \mathbf{a}^{k}}{\partial W^{rec}}$$
(18)

$$\frac{\partial^{+} a^{k}}{\partial W_{j}^{rec}} = \begin{bmatrix} \phi_{j}^{k} & 0 & \cdots & \cdots & 0 \\ 0 & \phi_{j}^{k} & \cdots & \cdots & 0 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & \cdots & \cdots & \cdots & \phi_{i}^{k} \end{bmatrix}$$
(19)

$$\frac{\partial \mathbf{a}^{t}}{\partial \mathbf{a}^{k}} = \frac{\partial \mathbf{a}^{t}}{\partial \mathbf{a}^{k+1}} \cdot diag(\sigma'(\mathbf{a}^{k})) \cdot W^{rec}$$

$$= \prod_{k=1}^{k} diag(\sigma'(\mathbf{a}^{i})) \cdot W^{rec}.$$
(20)

$$= \prod_{i=t-1}^{\kappa} diag(\sigma'(\mathbf{a}^i)) \cdot W^{rec}. \tag{21}$$

The derivatives with respect to the other variables are computed in a similar fashion.

Vanishing gradient: a sufficient condition

$$\frac{\partial \mathbf{a}^{t}}{\partial \mathbf{a}^{k}} = \prod_{i=t-1}^{k} diag(\sigma'(\mathbf{a}^{i})) \cdot W^{rec}. \tag{22}$$

Taking the singular value decomposition of W^{rec} :

$$W^{rec} = S \cdot D \cdot V^T \tag{23}$$

where S, V^T are squared orthogonal matrices and $D \triangleq diag(\mu_1, \mu_2, ..., \mu_r)$ is the diagonal matrix containing the singular values of W^{rec} . Hence:

$$\frac{\partial \mathbf{a}^t}{\partial \mathbf{a}^k} = \prod_{i=t-1}^k diag(\sigma'(\mathbf{a}^i)) \cdot S \cdot D \cdot V^T$$
(24)

Since U and V are orthogonal matrix, hence

$$||U||_2 = ||V^T||_2 = 1,$$

and

$$\|diag(\lambda_1, \lambda_2, ..., \lambda_r)\|_2 = \lambda_{max},$$

we get

$$\left\| \frac{\partial \boldsymbol{a}^{t}}{\partial \boldsymbol{a}^{k}} \right\|_{2} = \left\| \left(\prod_{i=t-1}^{k} diag(\sigma'(\boldsymbol{a}^{i})) \cdot S \cdot D \cdot V^{T} \right) \right\|_{2}$$
 (25)

$$\leq (\sigma'_{max} \cdot \mu_{max})^{t-k-1} \tag{26}$$

Solutions

Existent solutions

- ► Long short-term memory (LSTM). Hochreiter, Schmidhuber (1997)[1]
 - ▶ the network structure is modified with specialized "memory cells"
 - a truncated version of back-propagation is employed (but works also without it)
- ► Hessian-Free optimization (HF). Martens (2010) [3]
 - a second order method
 - a "cheap" approximation of the Hessian is employed
 - the quadratic sub-problem is solved through conjugate gradient + structural damping
- ▶ Pascanu, Bengio (2013) [5]
 - a first order method
 - uses a penalty to deal with the vanishing gradient problem

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Understanding the gradient structure: unfolding

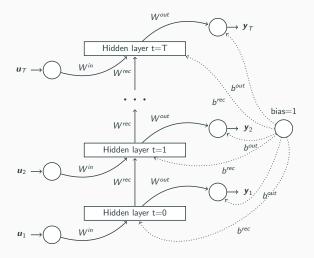


Figure 3: Unfolding of a RNN

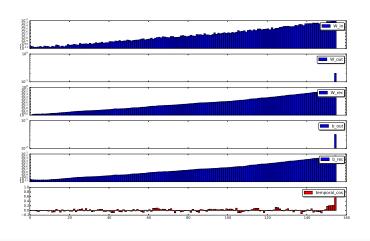
Understanding the gradient structure

It's easy to see that

$$\nabla L(\mathbf{x}) = \sum_{t=1}^{T} \nabla L_{|t}(\mathbf{x}), \tag{27}$$

where $\nabla L_{|t}$ is the gradient computed imagining to replicate all the variables for each time step and taking the derivatives w.r.t. the variables for step t

Temporal gradient norms: an illustration



A new proposal

The idea is to use the structure of the gradient to compute a "descent" direction which does not suffer from the vanishing problem.

normalize the temporal components:

$$g_t(\mathbf{x}) = \frac{\nabla L_{|t}(\mathbf{x})}{\|\nabla L_{|t}(\mathbf{x})\|}.$$
 (28)

combine the normalized gradients in a simplex:

$$g(\mathbf{x}) = \sum_{t=1}^{T} \beta_t \cdot g_t(\mathbf{x}). \tag{29}$$

with $\sum_{t=1}^{T} \beta_t = 1, \beta_t > 0$ (randomly picked at each iteration).

exploit the gradient norm:

$$d(\mathbf{x}) = \|\nabla L(\mathbf{x})\| \frac{g(\mathbf{x})}{\|g(\mathbf{x})\|}.$$
 (30)

Open issues

Open Issues: Initialization

- ► Some tasks, like the XOR one, are still "unresolved" (even for the other approaches). They cannot be solved with most of the seeds used for the initialization
- ▶ it seems to be an initialization matter

Popular strategies for initialization are:

- "small random weights", usually drawn from gaussian or uniform distribution with zero mean (Pascanu).
- sparse initialization: only some weights are actually sampled from a distribution, the other are set to zero (HF).
- ► ESN-like initialization[2].

Open Issues: Learning rate

- ▶ the **learning rate** is usually tuned by hand, there is no convergence theory for SGD in the non convex case
- ▶ a gradient clipping technique is often employed:

Algorithm 2: Gradient clipping

- 1 $\mathbf{g} \leftarrow \nabla_{\mathbf{x}} L$
- 2 if $\|\mathbf{g}\| \ge threshold$ then
- $oldsymbol{g} \leftarrow rac{ ext{threshold}}{\|oldsymbol{g}\|} oldsymbol{g}$
- 4 end
- some momentum or averaging technique often yield better convergence time, again tuned by hand

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