Welcome and Introduction

Philip Schulz and Wilker Aziz

https:
//github.com/philschulz/VITutorial

About us ...

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- Research associate at UvA
- VI, Sampling methods, Machine Translation

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- PhD candidate at UvA
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- ▶ VI, Machine Translation, Bayesian Models

Problems

Supervised problems: "learn a distribution over observed data"

sentences in natural language, images, videos, . . .

Unsupervised problems: "learn a distribution over observed and unobserved data"

sentences in natural language + parse trees, images + bounding boxes . . .

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▶ sentences, images, ...

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$$X \sim \mathsf{Cat}(\pi_1, \dots, \pi_K)$$
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and proceed to estimate parameters that assign maximum likelihood to observations

Multiple problems, same language



(Conditional) Density estimation

Side information (ϕ) **Parsing**

Observation (x)its syntactic/semantic a sentence

parse tree/graph

Translation a sentence its translation

Captioning an image caption in English

Entailment a text and hypothesis

entailment relation

Where does deep learning kick in?

Let ϕ be all side information available e.g. deterministic inputs/features

Have neural networks predict parameters of our probabilistic model

$$X|\phi \sim \mathsf{Cat}(\pi_{\theta}(\phi))$$
 or $X|\phi \sim \mathcal{N}(\mu_{\theta}(\phi), \sigma_{\theta}(\phi)^2)$

and proceed to estimate parameters w of the NNs

Task-driven feature extraction

Often our side information ϕ is itself some high dimensional data

- $ightharpoonup \phi$ is a sentence and x a tree
- $ightharpoonup \phi$ is the source sentence and x is the target
- lacktriangledown ϕ is an image and x is a caption

and part of the job of the NNs that parametrise our models is to also deterministically encode that input in a low-dimensional space

NN as efficient parametrisation

From the statistical point of view NNs do not generate data

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Prediction is done by a decision rule outside the statistical model

e.g. beam search

Let $p(x|\theta)$ be the probability of an observation x and θ refer to all of its parameters e.g. parameters of NNs involved

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$$\mathcal{L}(\theta|x^{(1:N)}) = \log \prod_{s=1}^{N} p(x^{(s)}|\theta)$$

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$$= \sum_{s=1}^{N} \log p(x^{(s)}|\theta)$$

MLE via gradient-based optimisation

If the log-likelihood is **differentiable** and **tractable** then backpropagation can give us the gradient

$$\mathbf{\nabla}_{ heta} \mathcal{L}(heta|x^{(1:N)}) = \mathbf{\nabla}_{ heta} \sum_{s=1}^{N} \log p(x^{(s)}| heta)$$

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and we can update θ in the direction

$$\gamma \nabla_{\theta} \mathcal{L}(\theta|x^{(1:N)})$$

to attain a local optimum of the likelihood function

$$abla_{ heta} \mathcal{L}(heta|x^{(1:N)}) = \underbrace{\sum_{s=1}^{N}
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For large N, computing the gradient is inconvenient

$$\begin{split} \boldsymbol{\nabla}_{\boldsymbol{\theta}} \mathcal{L}(\boldsymbol{\theta} | \boldsymbol{x}^{(1:N)}) &= \sum_{s=1}^{N} \boldsymbol{\nabla}_{\boldsymbol{\theta}} \log p(\boldsymbol{x}^{(s)} | \boldsymbol{\theta}) \\ &= \sum_{s=1}^{N} \frac{1}{N} N \boldsymbol{\nabla}_{\boldsymbol{\theta}} \log p(\boldsymbol{x}^{(s)} | \boldsymbol{\theta}) \\ &= \sum_{s=1}^{N} \mathcal{U}(\boldsymbol{s} | \boldsymbol{1} / \boldsymbol{N}) N \boldsymbol{\nabla}_{\boldsymbol{\theta}} \log p(\boldsymbol{x}^{(s)} | \boldsymbol{\theta}) \\ &= \mathbb{E}_{\boldsymbol{S} \sim \mathcal{U}(\boldsymbol{1} / \boldsymbol{N})} \left[N \boldsymbol{\nabla}_{\boldsymbol{\theta}} \log p(\boldsymbol{x}^{(s)} | \boldsymbol{\theta}) \right] \end{split}$$

S selects data points uniformly at random

Stochastic optimisation

For large N, we can use a gradient estimate

$$\nabla_{\theta} \mathcal{L}(\theta|x^{(1:N)}) = \underbrace{\mathbb{E}_{S \sim \mathcal{U}(^{1}/N)} \left[N \nabla_{\theta} \log p(x^{(S)}|\theta) \right]}_{\text{expected gradient :)}}$$

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and take a step in the direction

$$\gamma \frac{N}{M} \nabla_{\theta} \mathcal{L}(\theta | x^{(s_1:s_M)})$$

where $x^{(s_1:s_M)}$ is a random mini-batch of size M

DL in NLP recipe

Maximum likelihood estimation

 tells you which loss to optimise (i.e. negative log-likelihood)

Automatic differentiation (backprop)

"give me a tractable forward pass and I will give you gradients"

Stochastic optimisation powered by backprop

general purpose gradient-based optimisers

Tractability is central

Likelihood gives us a differentiable objective to optimise for

but we need to stick with tractable likelihood functions

When do we have intractable likelihood?

Unsupervised problems contain unobserved random variables

$$p(x, z | \theta) = \overbrace{p(z)}^{\text{prior}} \underbrace{p(x | z, \theta)}_{\text{observation model}}$$

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Unsupervised problems contain unobserved random variables

$$p(x, z|\theta) = \overbrace{p(z)}^{\text{prior}} \underbrace{p(x|z, \theta)}_{\text{observation model}}$$

thus assessing the marginal likelihood requires marginalisation of latent variables

$$p(x|\theta) = \int p(x,z|\theta) dz = \int p(z)p(x|z,\theta) dz$$

Examples of latent variable models

Discrete latent variable, continuous observation

$$p(x|\theta) = \underbrace{\sum_{c=1}^{K} \mathsf{Cat}(c|\pi_1, \dots, \pi_K) \underbrace{\mathcal{N}(x|\mu_{\theta}(c), \sigma_{\theta}(c)^2)}_{\mathsf{forward passes}}}_{\mathsf{too many forward passes}}$$

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Continuous latent variable, discrete observation

$$p(x|\theta) = \underbrace{\int \mathcal{N}(z|0,I) \underbrace{\operatorname{Cat}(x|\pi_{\theta}(z))}_{\text{forward pass}} dz}_{\text{infinitely many forward passes}}$$

Some reasons

 organise a massive collection of data e.g. LDA

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- organise a massive collection of data e.g. LDA
- learn from unlabelled data e.g. semi-supervised learning
- learn from little data e.g. Bayesian NNs
- induce discrete representations
 e.g. parse trees, dependency graphs,
 permutations, alignments

Probabilistic models parametrised by neural networks

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- explicit modelling assumptions one of the reasons why there's so much interest
- but requires efficient inference which is the reason why we are here today